

**2014 OU2 GROUNDWATER INVESTIGATION
VPB 154
BETHPAGE, NY**

Prepared for:



**Department of the Navy
Naval Facilities Engineering Command, Mid-Atlantic
9742 Maryland Ave.
Norfolk, VA 23511-3095**

**Comprehensive Long-Term Environmental Action Navy
Contract Number N62470-11-D-8013**

CTO WE15

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List of Acronyms and Abbreviations

| | |
|----------|---|
| AOC | Area of Concern |
| bgs | below ground surface |
| DoD | Department of Defense |
| ELAP | Environmental Laboratory Accreditation Program |
| EPA | Environmental Protection Agency, United States |
| ft | feet |
| GOCO | Government-Owned Contractor-Operated |
| GPS | Global Positioning System |
| IDW | Investigation Derived Waste |
| IR | Installation Restoration |
| Katahdin | Katahdin Analytical Services, Inc |
| NAD | North American Datum |
| NAVD | North American Vertical Datum |
| NAVFAC | Naval Facilities Engineering Command |
| NG | Northrop Grumman |
| NWIRP | Naval Weapons Industrial Reserve Plant |
| NYSDEC | New York State Department of Environmental Conservation |
| OU | Operable Unit |
| PCBs | Polychlorinated Biphenyls |
| PCE | Tetrachloroethene |
| PID | Photoionization Detector |
| POTW | Publicly Owned Treatment Works |
| PPE | Personal Protective Equipment |
| SAP | Sampling and analysis plan |
| SVOC | Semivolatile Organic Compounds |
| TCE | Trichloroethene |
| TCL | Target Compound List |
| TCLP | Toxicity Characteristic Leaching Procedure |
| TOC | Total Organic Carbon |
| UFP | United Federal Programs |
| VOC | Volatile Organic Compounds |
| VPB | Vertical Profile Boring |

1.0 PROJECT BACKGROUND

Resolution Consultants has prepared this Data Summary Report for the Naval Facilities Engineering Command (NAVFAC), Mid-Atlantic under contract task order WE15 Contract N62470-11-D-8013. This report describes vertical profile boring (VPB) installation activities (specifically at the VPB 154 location) in 2014 for the Naval Weapons Industrial Reserve Plant (NWIRP) Bethpage Operable Unit (OU) 2 Site 1 offsite plume. NWIRP Bethpage is located in east-central Nassau County, Long Island, New York, approximately 30 miles east of New York City (Figure 1).

1.1 Scope and Objectives

This data summary report provides information on the installation of VPB 154. The purpose of the VPB 154 investigation was to determine subsurface conditions and contaminant levels in the southwestern portion of the plume and define the western and southern extent of the 108 Hot Spot (defined as an area >1000ppb of total volatile organic compounds [VOCs] north of Hempstead Turnpike). VPB locations within the general vicinity of VPB 154 are shown in Figure 2. VPB 154 was completed to 950 feet (ft) below ground surface (bgs).

Field tasks were conducted in 2014 in accordance with the *United Federal Programs Sampling and Analysis Plan (UFP SAP)*, Bethpage, New York and the UFP SAP Addendum Installation of Vertical Profile Borings and Monitoring Wells (Resolution Consultants, 2013). The field investigation included completing one vertical profile boring, groundwater grab samples, geophysical logging and surveying.

Documentation of these activities is included in Appendix A of this report.

1.2 Site History

NWIRP Bethpage is in the Hamlet of Bethpage, Town of Oyster Bay, New York. Since its inception in 1941, the plant's primary mission was the research prototyping, testing, design, engineering, fabrication, and primary assembly of military aircraft. The facilities at NWIRP included four plants used for assembly and prototype testing, a group of quality control laboratories, two warehouse complexes (north and south), a salvage storage area, water recharge basins, the Industrial Wastewater Treatment Plant, and several smaller support buildings.

The Navy's property originally totaled 109.5 acres and was formerly a Government-Owned Contractor-Operated (GOCO) facility that was operated by Northrop Grumman (NG) until September 1998. Prior to 2002, the NWIRP property was bordered on the north, west, and south

by current or former NG facilities, and on the east by a residential neighborhood. By March 2008, approximately 100 acres of NWIRP property were transferred to Nassau County in three separate actions. The remaining 9 acres and access easements were retained by the Navy to continue remedial efforts at Installation Restoration (IR) Site 1 – Former Drum Marshalling Area and Site 4 – Former Underground Storage Tanks (Area of Concern [AOC] 22). A parcel of land connecting the two sites was also retained. Currently, the 9-acre parcel of NWIRP is bordered on the east by the residential neighborhood and on the north, south, and west by Nassau County property. Access to the NWIRP is from South Oyster Bay Road.

1.3 Geology and Hydrogeology

Overburden at the site consists of well over 1,000 ft of Cretaceous deposits overlying crystalline bedrock of the Hartland Formation. Overburden is divided into four geologic units: the upper Pleistocene deposits, the Magothy Formation, the clay member of the Raritan Formation ("Raritan Clay") and the Lloyd Sand member of the Raritan Formation ("Lloyd Sand") (Geraghty and Miller, 1994).

The upper Pleistocene ranges in thickness from approximately 50 to 100 ft and consists of till and outwash deposits of medium to coarse sand and gravel with lenses of fine sand, silt and clay (Smolensky and Feldman, 1990); these deposits form the Upper Glacial Aquifer. Directly underlying this unit is the Magothy Formation with a thickness of 650 to 900 ft bgs observed onsite. The Magothy is characterized by fine to medium sands and silts interbedded with zones of clays, silty sands and sandy clays. Sand and gravel lenses are found in some areas between depths of 600 and 875 ft bgs; these deposits form the Magothy Aquifer.

Investigations performed by the Navy since 2012 indicate that the bottom of the Magothy (top of the Raritan Clay) can extend to depths of 700 to greater than 1,000 ft bgs. The top of the Raritan Clay deepens to the south southeast, as evidenced by clay depths of 1,000 ft bgs (or more) intexborings installed offsite. The Raritan Clay Unit is of continental origin and consists of clay, silty clay, clayey silt, and fine silty sand. This member acts as a confining layer over the Lloyd Sand Unit. The Lloyd Sand Unit is also of continental origin, having been deposited in a large fresh water lacustrine environment. The material consists of fine to coarse-grained sands, gravel, interbedded clay, and silty sand. These deposits form the Lloyd Aquifer.

The Upper Glacial Aquifer and the Magothy Aquifer comprise the aquifers of interest at the NWIRP. Regionally, these formations are generally considered to form a common, interconnected aquifer as

the coarse nature of each unit near their contact and the lack of any regionally confining clay unit allows for the unrestricted flow of groundwater between the formations.

The Magothy Aquifer is the major source of public water in Nassau County. The most productive water bearing zones are the discontinuous lenses of sand and gravel that occur within the siltier matrix. The major water-bearing zones are course sand and gravel lenses located in the lower portion of the Magothy. The Magothy Aquifer is commonly regarded to function overall as an unconfined aquifer at shallow depths and a confined aquifer at deeper depths. The drilling program at the NWIRP has revealed that clay zones beneath the facility are common but laterally discontinuous. No confining clay units of facility-wide extent have been encountered.

Groundwater is encountered at a depth of approximately 50 ft bgs at the facility. Historically, because of pumping and recharge at the facility, groundwater depths have been measured to range from 40 to 60 ft bgs. The groundwater flow in the area is to the south-southeast.

2.0 FIELD PROGRAM

Field investigation activities at VPB 154 consisted of drilling, sampling, soil/groundwater analysis, geophysical logging, and surveying. Drilling during this investigation was performed by Delta Well and Pump Company of Ronkonkoma, New York. A description of these tasks is provided below.

2.1 Vertical Profile Borings

One vertical profile boring (VPB 154) was completed during this field effort between July 15, 2014 and September 3, 2014. The total depth of VPB 154 was 950 ft. The location is shown in Figure 2 and details are summarized in Table 1.

2.1.1 Drilling

VPB 154 was installed by drilling a 8-inch diameter hole using mud rotary drilling techniques. Drilling mud consisted of potable water and polymer-free sodium bentonite or equivalent. Drilling mud was contained and re-circulated in baffled, high capacity mud tubs. A sand separator was used intermittently to remove fines from circulation.

2.1.2 Sampling

A total of seven split spoon samples were collected from ground surface to the bottom of the boring. A change in geology was observed by the field geologist at 938 ft bgs and three split spoon samples were subsequently collected to confirm the presence of the Raritan Clay. Samples were logged by the field geologist and screened for Volatile Organic Compounds (VOCs) utilizing a photoionization detector (PID). A detailed boring log for VPB 154 is included in Appendix A.

Groundwater grab samples were collected every 50 ft for the first 200 ft of borehole depth. After the first 200 ft, groundwater grab samples were collected approximately every 20 ft until the boring terminated in the Raritan. Groundwater grab samples were collected with a hydropunch sampler and analyzed for VOCs using Environmental Protection Agency (EPA) Method 8260C. The groundwater grab samples were analyzed by Katahdin Analytical Services (Katahdin), a Department of Defense (DoD), Environmental Laboratory Accreditation Program (ELAP), and New York State Department of Environmental Conservation (NYSDEC)-certified laboratory. During the collection of groundwater grab samples, field parameters were measured (pH, temperature, specific conductivity, oxidation reduction potential, dissolved oxygen, and turbidity). Data validation was performed by Resolution Consultants. Groundwater grab sample logs, data validation packages, and analytical data tables are included in Appendix A.

One soil sample was collected for laboratory analysis for total organic carbon (TOC) by EPA series SW-846 method 9060A. During drilling, air sampling was conducted under a Community Air Monitoring Plan. One air sample was collected per VPB using Summa canisters and submitted for laboratory analysis by EPA Method TO-15. All analyses were performed or sub-contracted by Katahdin. Data validation of both TOC and air data was performed by Resolution Consultants. Data validation packages and analytical data tables are included in Appendix A.

2.1.3 Geophysics

Borehole geophysical logs (gamma) were recorded after the borehole was drilled but prior to the removal of drill rods. A Mount Sopris Instrument model 2PGA-100 poly gamma was used. Starting at the top of the hole, the probe was advanced at a maximum rate of 12 ft per minute. A copy of the log was printed in the field for review once the probe reached the bottom of the borehole. The instrument was then raised to the top of the boring and a second log was generated and printed in the field. The down hole gamma log sheets and plots comparing the gamma log with trichloroethene (TCE) and tetrachloroethene (PCE) concentrations from hydropunch samples are included in Appendix A.

2.2 Decontamination and Investigation Derived Waste (IDW)

Resolution Consultants utilized dedicated and disposable sampling equipment when possible to avoid the potential for cross-contamination of samples. The sampling equipment included dedicated plastic scoops, disposable Teflon or polyethylene tubing, disposable gloves, and laboratory supplied sample bottles. Hand held equipment, split spoons, and the hydropunch were decontaminated using Liquinox and water wash, a potable water rinse, followed by a distilled water rinse. Water was collected in 5-gallon pails or 55-gallon drums.

As part of the IDW management practices and in accordance with the SAP, the investigation waste (consisting of soil cuttings, drilling muds, IDW fluids, and personal protective equipment (PPE)) generated during the boring installation was containerized and staged at NWIRP Bethpage. IDW solids were characterized and disposed of properly. Representative samples from each roll off were submitted to Katahdin for analysis of:

- Target Compound List (TCL) VOCs
- TCL Semi-volatile Organic Compounds (SVOCs)
- Toxicity Characteristic Leaching Procedure (TCLP) Metals

-
- Polychlorinated Biphenyls (PCBs)
 - Total petroleum hydrocarbons
 - Corrosivity
 - Ignitability
 - Reactive Cyanide
 - Reactive Sulfide
 - Paint Filter

IDW water was containerized in frac tanks and stored at NWIRP Bethpage for characterization and ultimate disposal to the Publicly Owned Treatment Works (POTW), in accordance with the facilities existing discharge permit. A representative water sample was collected from each frac tank and submitted to Katahdin for analysis of VOCs via Method SW 624, pH via Method SW 9040B, PCBs via Method 8082 and Total Metals via Method SW 846 (all waters). To the extent feasible, soil and water were not mixed. All analytical criteria were met for disposal of soil and water.

2.3 Surveying

A survey of the boring location was conducted at the end of fieldwork by C. T. Male, Inc., of Latham, NY, under the direct supervision of Resolution Consultants. The location was tied into the existing base map developed for this investigation. The survey elevation is referenced to the North American Vertical Datum (NAVD) 1988 and has a vertical accuracy of 0.01 foot. Vertical control is based on observations of COR Stations Queens and Central Islip. The horizontal location is referenced to the North American Datum (NAD) 1983 (2011) N.Y. Long Island Zone 3104 and has an accuracy of 0.1 foot. Local horizontal and vertical control is based on Global Positioning System (GPS) observations using the NYS Net Real Time Network.

A table of survey data (ground, latitude/longitude and northing/easting) and a survey map is included in Appendix A.

3.0 REFERENCES

Geraghty and Miller, Inc., 1994. *Remedial Investigation Report, Grumman Aerospace Corporation, Bethpage, New York*. Revised September 1994.

Naval Facilities Engineering Command (NAVFAC), 2003. *Record of Decision Naval Weapons Industrial Reserve Plant Bethpage, New York, Operable Unit 2 – Groundwater*, NYS Registry: 1-30-003B. April.

Resolution Consultants, 2013. *United Federal Programs Sampling and Analysis Plan, Site OU-2 Offsite TCE Groundwater Plume Investigation*, Bethpage, New York. April.

Resolution Consultants, 2013. UFP SAP Addendum, *Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol*. November.

Resolution Consultants, 2013. UFP SAP Addendum, *Installation of Vertical Profile Borings and Monitoring Wells*. December.

Smolensky, D., and Feldman, S., 1990. *Geohydrology of the Bethpage-Hicksville-Levittown Area, Long Island, New York*, U.S. Geological Survey Water-Resourced Investigations Report 88-4135, 25 pp.

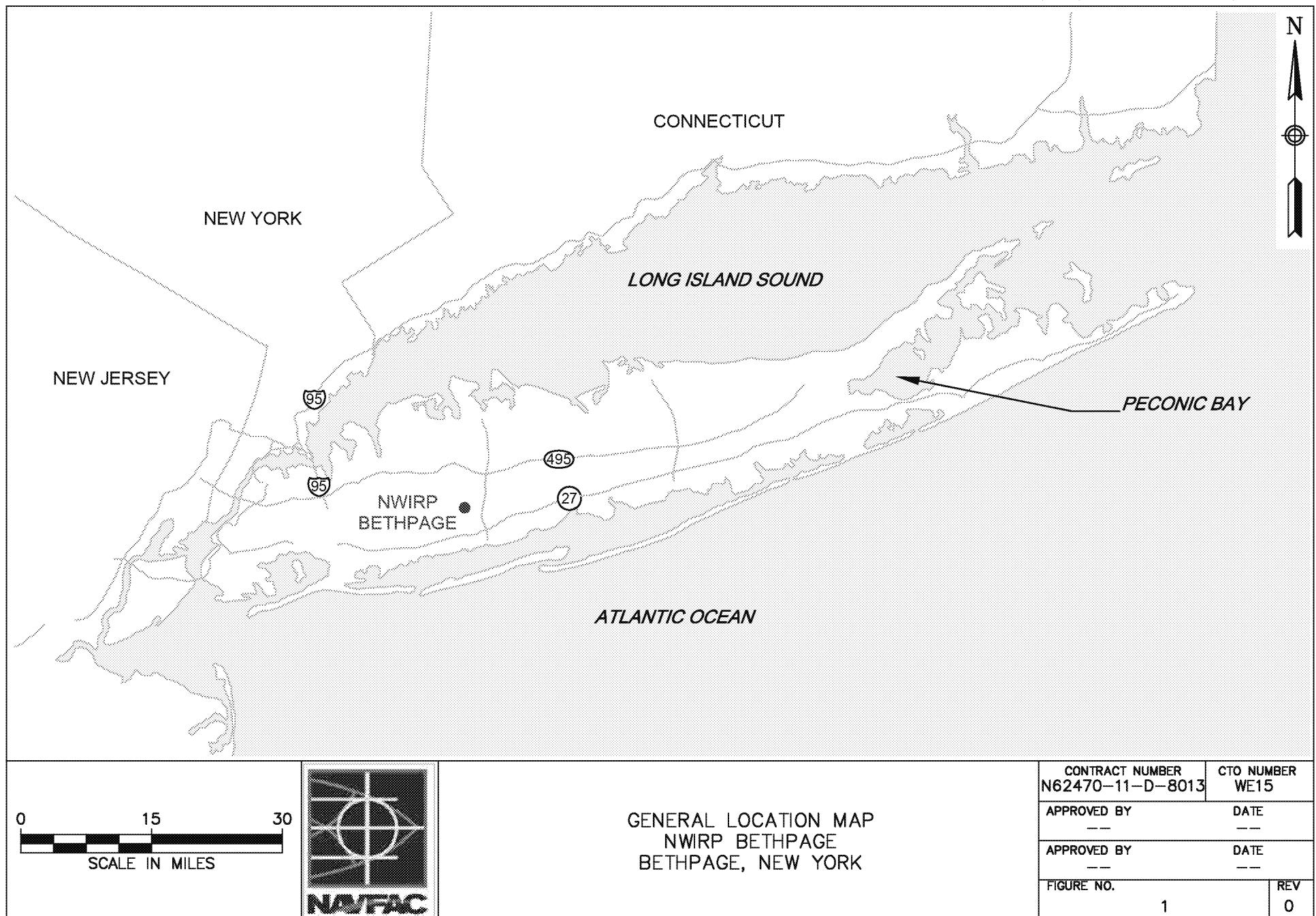
Tables

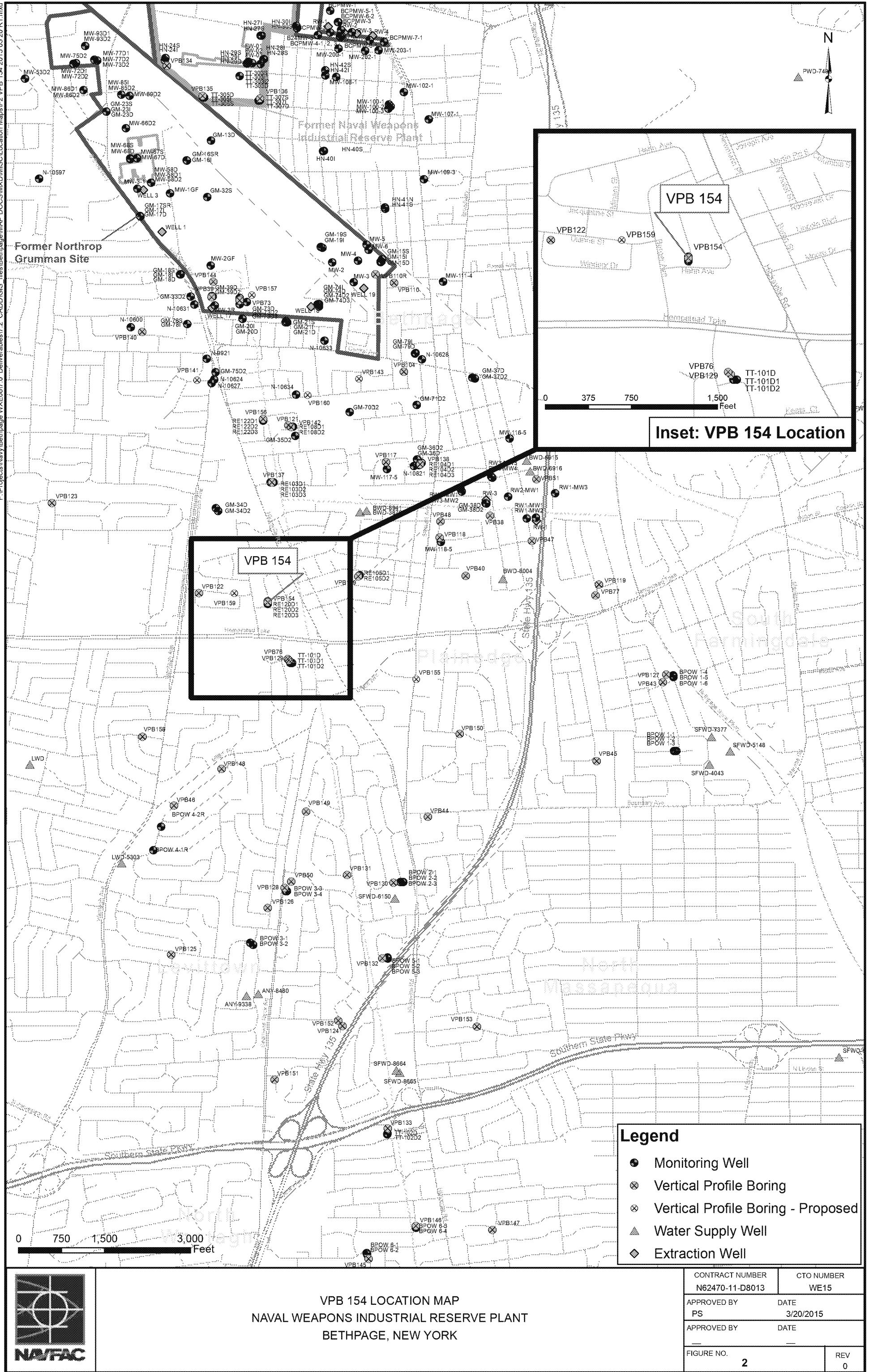
TABLE 1
VERTICAL PROFILE BORING SUMMARY
2014 OU2 GROUNDWATER INVESTIGATION
NWIRP BETHPAGE, NY

| BORING | BORING START DATE | BORING COMPLETION DATE | GROUND ELEVATION (MSL) | TOTAL DEPTH (ft bgs) | SURFACE CASING SET AT (ft bgs) | NO. OF SPOON SAMPLES | GAMMA LOG (ft bgs) | NO. GW SAMPLES COLLECTED/ATTEMPTED* | TOC SAMPLES | DATE OF AIR SAMPLE | MONITORING WELLS INSTALLED AT LOCATION |
|---------|-------------------|------------------------|------------------------|----------------------|--------------------------------|----------------------|--------------------|-------------------------------------|------------------|--------------------|--|
| VPB 154 | 7/15/2014 | 9/3/2014 | 85.86 | 950 | 53 | 7 | 947 | 41/49 | 422 - 424 ft bgs | 9/2/2014 | Pending |

* includes field duplicates

Figures





Appendix A

VPB 154

Section 1
VPB 154 Boring and Gamma Logs

Resolution Consultants

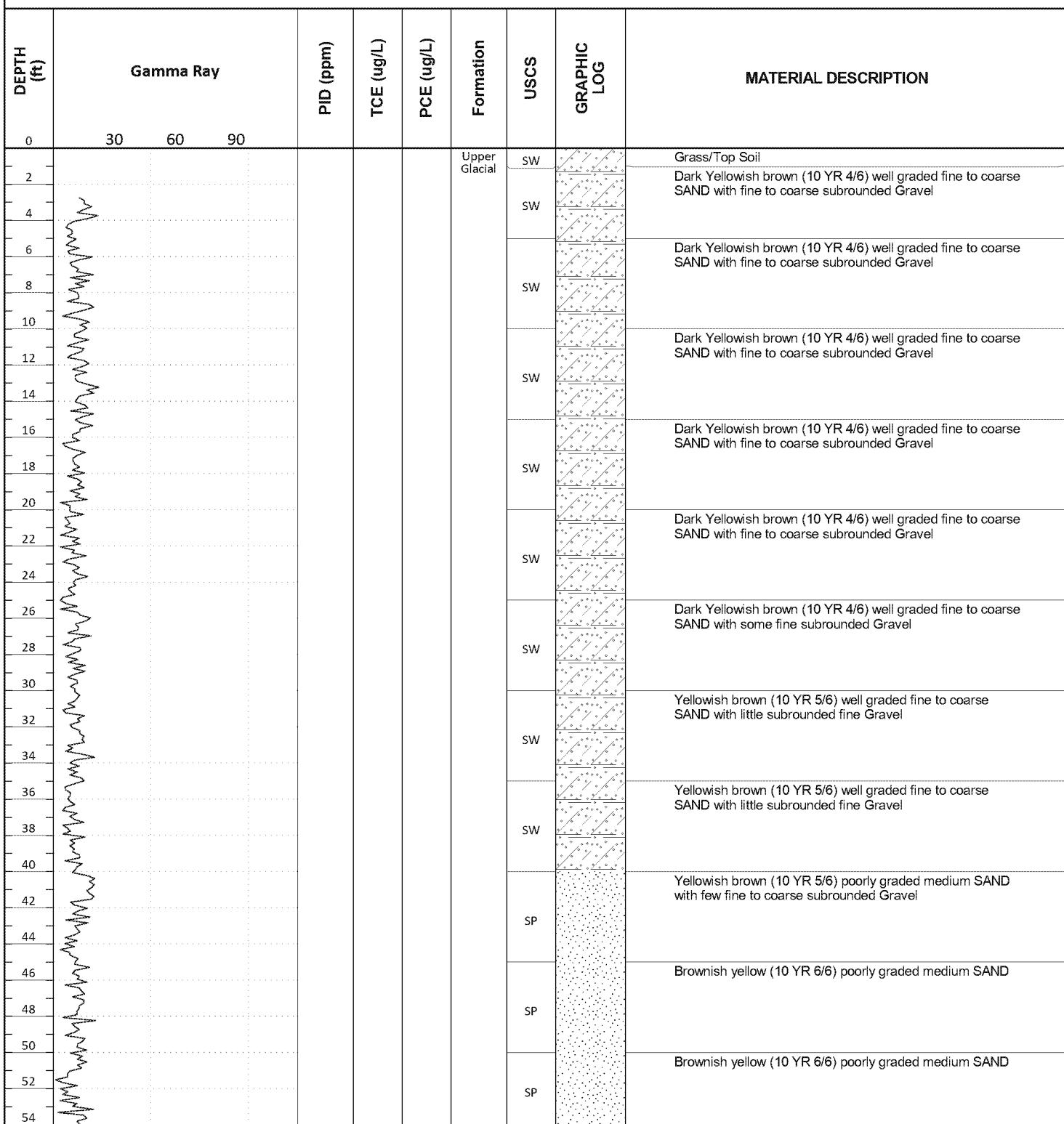
Boring Log

BORING #: VPB154

Sheet 1 of 16

| | | | | | | |
|--|--|--|--|--|-------------------------------------|--------------------------|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | | | | | Logged By: V. Varricchio |
| Location: Shelly Dr. and Hahn Ave., Bethpage, NY | Northing: 204605.97 Easting: 1125061.14 | | | | Drilling Company: Delta Well & Pump | |
| Project #: 60266526 | Ground Elevation (ft amsl): 85.86 | | | | Well Screen Interval (ft): NA | |
| Start Date: 7/15/2014 | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | | | | Water Level (ft): NA | |
| Finish Date: 9/3/2014 | | | | | Total Depth (ft): 950.0 | |

Mud Rotary Drilling Note: Unless denoted by a splitspoon sample (indicated by the presence of a PID reading), boundaries between strata are approximate only and may be transitional because they are based on screened wash samples collected during mud rotary drilling at 5 ft. intervals.



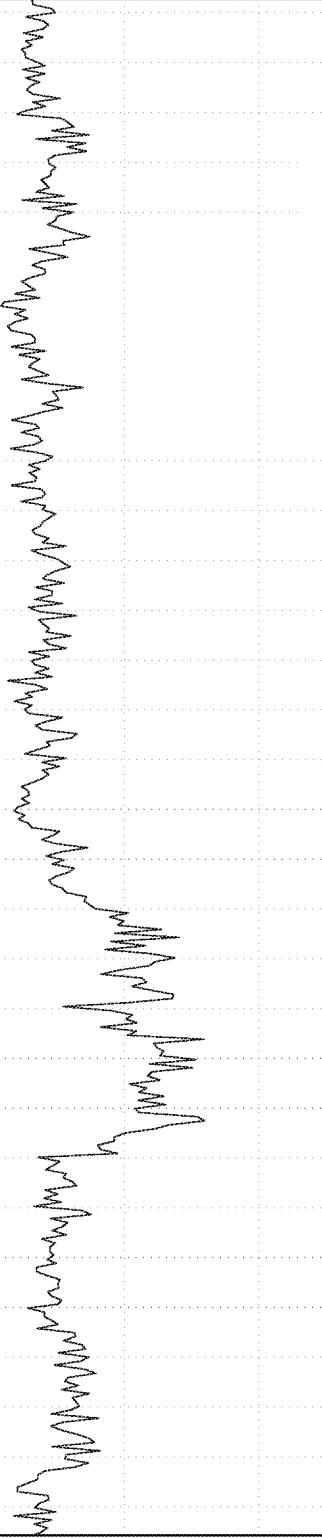
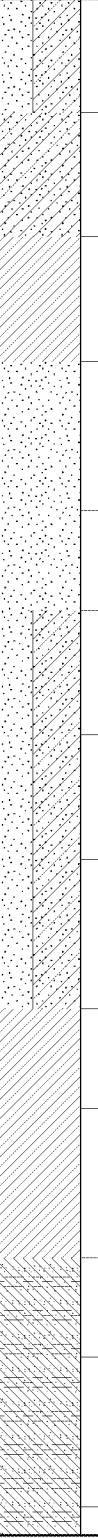
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| DEPTH (ft) | Gamma Ray | | | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|----|----|-----------|------------|------------|------------------|-------|----------------|---|
| 54 | 30 | 60 | 90 | | | | | | | |
| 56 | | | | | | | Upper Glacial | SP | | Brownish yellow (10 YR 6/6) well graded medium to coarse SAND with some Silt and fine subrounded gravel |
| 58 | | | | | | | | SP-SM | | |
| 60 | | | | < 0.5 | < 0.5 | | | SW | | Brownish yellow (10 YR 6/6) well graded fine to coarse SAND with fine to coarse subrounded Gravel |
| 62 | | | | | | | | SP | | Reddish yellow (7.5 YR 6/6) poorly graded medium SAND with trace fine subrounded Gravel |
| 64 | | | | | | | | SP | | Reddish yellow (7.5 YR 6/6) poorly graded medium SAND with trace fine subrounded Gravel |
| 66 | | | | | | | | SP | | Reddish yellow (5 YR 6/8) poorly graded medium SAND with fine subrounded Gravel |
| 68 | | | | | | | | SP | | Very pale brown (10 YR 7/3) poorly graded medium SAND |
| 70 | | | | | | | | SP | | |
| 72 | | | | | | | | SW | | Reddish yellow (7.5 YR 6/6) well graded fine to coarse SAND with trace lean Clay and subrounded fine gravel |
| 74 | | | | | | | | GW | | Very pale brown (10 YR 7/4) well graded subrounded GRAVEL with medium to coarse Sand, trace lean clay |
| 76 | | | | | | | | GW | | |
| 78 | | | | | | | | SW-GW | | Very pale brown (10 YR 7/4) well graded subrounded GRAVEL with medium to coarse Sand, trace lean clay |
| 80 | | | | | | | | SW-GW | | Light yellowish brown (10 YR 6/4) Gravely well graded medium to coarse SAND |
| 82 | | | | | | | | SW-GW | | Light yellowish brown (10 YR 6/4) Gravely well graded medium to coarse SAND |
| 84 | | | | | | | | GP | | Light yellowish brown (10 YR 6/4) poorly graded subrounded fine GRAVEL with few medium to coarse Sand |
| 86 | | | | | | | | | | |
| 88 | | | | | | | | | | |
| 90 | | | | | | | | | | |
| 92 | | | | | | | | | | |
| 94 | | | | | | | | | | |
| 96 | | | | | | | | | | |
| 98 | | | | | | | | | | |
| 100 | | | | < 0.5 | < 0.5 | | Magothy | | | |
| 102 | | | | | | | | | | |
| 104 | | | | | | | | | | |
| 106 | | | | | | | | | | |
| 108 | | | | | | | | | | |
| 110 | | | | | | | | | | |
| 112 | | | | | | | | | | |
| 114 | | | | | | | | | | |

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| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|--|
| 116 | 30 60 90 | | | | | | | |
| 118 | | | | | Magothy | GP | | Light yellowish brown (10 YR 6/4) well graded subrounded GRAVEL with trace medium to coarse Sand |
| 120 | | | | | | GW | | |
| 122 | | | | | | GW-SW | | Light brown (7.5 YR 6/4) Sandy well graded subrounded GRAVEL with trace lean clay |
| 124 | | | | | | CL | | Brown (10 YR 5/3) Sandy lean CLAY with trace fine subrounded gravel |
| 126 | | | | | | GC | | Pale brown (10 YR 6/3) Clayey coarse subrounded GRAVEL with few well graded medium coarse sand |
| 128 | | | | | | SP | | Pale brown (10 YR 6/3) poorly graded fine SAND |
| 130 | | | | | | SP | | Pale brown (10 YR 6/3) poorly graded fine SAND with trace fine subrounded Gravel and lean clay |
| 132 | | | | | | ML | | Light yellowish brown (10 YR 6/4) fine to medium Sandy SILT with trace iron nodules |
| 134 | | | | | | ML | | Light yellowish brown (10 YR 6/4) fine to medium Sandy SILT with trace iron nodules |
| 136 | | | | | | SP | | Light brownish gray (10 YR 6/2) poorly graded medium SAND with trace iron nodules and Clay |
| 138 | | | | | | SP | | Light brownish gray (10 YR 6/2) poorly graded medium SAND with trace iron nodules and Clay |
| 140 | | | | | | SP | | Pale brown (2.5 Y 8/2) poorly graded medium SAND with trace iron nodules |
| 142 | | | | | | SP | | Pale brown (2.5 Y 8/2) poorly graded medium SAND with trace iron nodules |
| 144 | | | | | | | | |
| 146 | | | | | | | | |
| 148 | | | | | | | | |
| 150 | | < 0.5 | | < 0.5 | | | | |
| 152 | | | | | | | | |
| 154 | | | | | | | | |
| 156 | | | | | | | | |
| 158 | | | | | | | | |
| 160 | | | | | | | | |
| 162 | | | | | | | | |
| 164 | | | | | | | | |
| 166 | | | | | | | | |
| 168 | | | | | | | | |
| 170 | | | | | | | | |
| 172 | | | | | | | | |
| 174 | | | | | | | | |
| 176 | | | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | | | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|--|----|----|-----------|------------|------------|-----------|-------|---|--|
| | 30 | 60 | 90 | | | | | | | |
| 178 |  | | | | | | Magothy | SP-SC |  | Pale brown (10 YR 6/5) poorly graded medium SAND with little Clay and trace iron nodules (continued) |
| 180 | | | | | | | | SC | | |
| 182 | | | | | | | | CL | | |
| 184 | | | | | | | | SP | | |
| 186 | | | | | | | | SP | | |
| 188 | | | | | | | | SP-SC | | |
| 190 | | | | | | | | SP-SC | | |
| 192 | | | | | | | | SP-SC | | |
| 194 | | | | | | | | SP-SC | | |
| 196 | | | | | | | | SP-SC | | |
| 198 | | | | | | | | SP-SC | | |
| 200 | | | | 31 | 3.2 | | | SP-SC | | |
| 202 | | | | | | | | SP-SC | | |
| 204 | | | | | | | | SP-SC | | |
| 206 | | | | | | | | SP-SC | | |
| 208 | | | | | | | | SP-SC | | |
| 210 | | | | | | | | SP-SC | | |
| 212 | | | | | | | | SP-SC | | |
| 214 | | | | | | | | SP-SC | | |
| 216 | | | | | | | | SP-SC | | |
| 218 | | | | | | | | SP-SC | | |
| 220 | | | | | | | | SP-SC | | |
| 222 | | | | | | | | SP-SC | | |
| 224 | | | | | | | | SP-SC | | |
| 226 | | | | | | | | SP-SC | | |
| 228 | | | | | | | | SP-SC | | |
| 230 | | | | 180 | 4.8 | | | ML-CL |  | Dark gray (10 YR 4/1) lean CLAY with some poorly graded medium Sand, trace iron nodules |
| 232 | | | | | | | | ML-CL | | |
| 234 | | | | | | | | ML-CL | | |
| 236 | | | | | | | | ML-CL | | |
| 238 | | | | 190 | 5.1 | | | ML-CL | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | | |
|---------------|-----------|-----------|------------|------------|-----------|---------|-------------|----------------------|--|----|
| | | | | | | | | 30 | 60 | 90 |
| 240 | | | | 190 | 5.1 | Magothy | ML-CL | | Yellowish brown (10 YR 5/4) Clayey SILT with some medium Sand and trace iron nodules (continued) | |
| 242 | | | | | | | | | | |
| 244 | | | | | | | | | | |
| 246 | | | | | | | | | | |
| 248 | | | | | | | | | | |
| 250 | | | | | | | | | | |
| 252 | | | | | | | | | | |
| 254 | | | | | | | | | | |
| 256 | | | | | | | | | | |
| 258 | | | | | | | | | | |
| 260 | | | < 0.5 | < 0.5 | | | | | | |
| 262 | | | | | | | | | | |
| 264 | | | | | | | | | | |
| 266 | | | | | | | | | | |
| 268 | | | | | | | | | | |
| 270 | | | | | | | | | | |
| 272 | | | | | | | | | | |
| 274 | | | | | | | | | | |
| 276 | | | | | | | | | | |
| 278 | | | | | | | | | | |
| 280 | | | 200 | 7.1 | | | | | | |
| 282 | | | | | | | | | | |
| 284 | | | | | | | | | | |
| 286 | | | | | | | | | | |
| 288 | | | | | | | | | | |
| 290 | | | | | | | | | | |
| 292 | | | | | | | | | | |
| 294 | | | | | | | | | | |
| 296 | | | | | | | | | | |
| 298 | | | | | | | | | | |
| 300 | | | 200 | 8.0 | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|---|
| | | | | | | | | 30 |
| 302 | | | | | Magothy | SP | | Yellowish brown (10 YR 5/4) poorly graded medium SAND with little iron nodules and trace Clay |
| 304 | | | | | | SP | | Yellowish brown (10 YR 5/4) poorly graded fine SAND with trace iron nodules |
| 306 | | | | | | SP | | Yellowish brown (10 YR 5/4) well graded fine to medium SAND with iron nodules and trace Clay |
| 308 | | | | | | SW | | Dark grayish brown (10 YR 4/2) lean CLAY with trace fine Sand and iron nodules |
| 310 | | | | | | CL | | Dark grayish brown (10 YR 4/2) lean CLAY with little fine Sand |
| 312 | | | | | | CL | | Grayish brown (10 YR 5/2) fine Sandy lean CLAY |
| 314 | | | | | | CL-ML | | Grayish brown (10 YR 5/2) Silty CLAY with little fine Sand and trace iron nodules |
| 316 | | | | | | SP | | Grayish brown (10 YR 5/2) poorly graded fine SAND, trace iron nodules |
| 318 | | | | | | SP | | Yellowish brown (10 YR 5/4) poorly graded fine SAND, trace iron nodules |
| 320 | | | | | | SP | | Very dark gray (10 YR 3/1) Silty CLAY with little fine to medium Sand and iron nodules |
| 322 | | | | | | SP | | Very dark gray (10 YR 3/1) fat CLAY |
| 324 | | | | | | CH | | |
| 326 | | | | | | | | |
| 328 | | | | | | | | |
| 330 | | | | | | | | |
| 332 | | | | | | | | |
| 334 | | | | | | | | |
| 336 | | | | | | | | |
| 338 | | | | | | | | |
| 340 | | | | | | | | |
| 342 | | | | | | | | |
| 344 | | | | | | | | |
| 346 | | | | | | | | |
| 348 | | | | | | | | |
| 350 | | | | | | | | |
| 352 | | | | | | | | |
| 354 | | | | | | | | |
| 356 | | | | | | | | |
| 358 | | | | | | | | |
| 360 | | | | | | | | |
| 362 | | | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|---|
| 364 | | | | | Magothy | CH | | Very dark gray (10 YR 3/1) fat CLAY (continued) |
| 366 | | | | | | CH | | |
| 368 | | | | | | CH | | Very dark gray (10 YR 3/1) fat CLAY |
| 370 | | | | | | CH | | |
| 372 | | | | | | CH | | Very dark gray (10 YR 3/1) fat CLAY |
| 374 | | | | | | CH | | |
| 376 | | | | | | CH | | |
| 378 | | | | | | CH | | Black (10 YR 2/1) fat CLAY |
| 380 | | | | | | CH | | |
| 382 | | | | | | CH | | Very dark gray (10 YR 3/1) fat CLAY, trace Silt |
| 384 | | | | | | CH | | |
| 386 | | | | | | ML-CH | | |
| 388 | | | | | | ML-CH | | Dark gray (10 YR 4/1) Clayey SILT with few poorly graded fine Sand |
| 390 | | | | | | ML-CH | | |
| 392 | | | | | | ML-CH | | Dark gray (10 YR 4/1) Clayey SILT with little poorly graded fine Sand |
| 394 | | | | | | SM-SC | | |
| 396 | | | | | | SM-SC | | Dark gray (10 YR 4/1) Silty Clayey poorly graded fine SAND |
| 398 | | | | | | SM-SC | | |
| 400 | | | | | | SM-SC | | Dark gray (10 YR 4/1) Silty Clayey poorly graded fine SAND |
| 402 | | | | | | SM-SC | | |
| 404 | | | | | | SM-SC | | Dark gray (10 YR 4/1) Silty Clayey poorly graded fine SAND |
| 406 | | | | | | SM-SC | | |
| 408 | | | | | | SM-SC | | Dark gray (10 YR 4/1) Silty Clayey poorly graded fine SAND |
| 410 | | | | | | SP | | |
| 412 | | | | | | SP | | Grayish brown (10 YR 5/2) poorly graded fine SAND |
| 414 | | | | | | SP | | |
| 416 | | | | | | SP | | Gray (10 YR 6/1) poorly graded fine SAND |
| 418 | | | | | | SP | | |
| 420 | | | | | | SP | | Gray (10 YR 6/1) poorly graded fine SAND |
| 422 | | | | | | SP | | |
| 424 | | 0 | 170 | 4.8 | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | |
|---------------|-----------|-----------|------------|------------|-----------|-------|-------------|---|----|
| | | | | | | | | 30 | 60 |
| 426 | | | | | Magothy | SP | | Gray (10 YR 6/1) poorly graded fine SAND (continued) | |
| 428 | | | | | | SP-SM | | Dark gray (10 YR 4/1) poorly graded fine SAND with few Silt | |
| 430 | | | | | | CL | | Dark gray (10 YR 4/1) Sandy lean CLAY | |
| 432 | | | | | | CL | | Dark gray (10 YR 4/1) Sandy lean CLAY | |
| 434 | | | | | | CH | | Grayish brown (10 YR 5/2) fat CLAY | |
| 436 | | | | | | CH | | Gray (10 YR 5/1) fat CLAY, trace iron nodules | |
| 438 | | | | | | CH | | Grayish brown (10 YR 5/2) fat CLAY with little poorly graded fine Sand | |
| 440 | | 190 | 5.0 | | | SP | | Light gray (10 YR 7/1) poorly graded medium SAND, trace Clay | |
| 442 | | | | | | CL | | Grayish brown (10 YR 5/2) lean CLAY with some poorly graded medium Sand, trace iron nodules | |
| 444 | | | | | | SW | | Light grayish brown (10 YR 6/2) well graded fine to coarse SAND with trace Clay | |
| 446 | | | | | | SW | | Light grayish brown (10 YR 6/2) well graded fine to coarse SAND with trace Clay | |
| 448 | | | | | | SP-SC | | Light gray (10 YR 7/2) poorly graded medium SAND with few Clay, trace iron nodules | |
| 450 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND, trace iron nodules | |
| 452 | | | | | | | | | |
| 454 | | | | | | | | | |
| 456 | | | | | | | | | |
| 458 | | | | | | | | | |
| 460 | | 180 | 5.4 | | | | | | |
| 462 | | | | | | | | | |
| 464 | | | | | | | | | |
| 466 | | | | | | | | | |
| 468 | | | | | | | | | |
| 470 | | | | | | | | | |
| 472 | | | | | | | | | |
| 474 | | | | | | | | | |
| 476 | | | | | | | | | |
| 478 | | | | | | | | | |
| 480 | | | | | | | | | |
| 482 | | | | | | | | | |
| 484 | | | | | | | | | |
| 486 | | 160 | 6.6 | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|--|
| 486 | 30 60 90 | | | | | | | |
| 488 | | | | | Magothy | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND, trace iron nodules |
| 490 | | | | | | SC | | |
| 492 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND, trace iron nodules |
| 494 | | | | | | SC | | |
| 496 | | | | | | SW-SC | | |
| 498 | | | | | | SC | | Light brownish gray (10 YR 6/2) well graded fine to medium SAND with little Clay |
| 500 | | | | | | SC | | Grayish brown (10 YR 5/2) Clayey well graded fine to medium SAND, trace iron nodules |
| 502 | | | | | | CH | | Gray (10 YR 5/1) fat CLAY with little poorly graded fine Sand, trace iron nodules |
| 504 | | 160 | < 0.5 | | | SC | | Grayish brown (10 YR 5/2) Clayey well graded fine to medium SAND |
| 506 | | | | | | SC | | |
| 508 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND |
| 510 | | | | | | SC | | |
| 512 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND |
| 514 | | | | | | SC | | |
| 516 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine to medium SAND |
| 518 | | | | | | SC | | |
| 520 | | 340 | 3.4 | | | SC | | Grayish brown (10 YR 5/2) well graded fine to medium SAND with few Clay |
| 522 | | | | | | SC | | |
| 524 | | | | | | SC | | Dark grayish brown (10 YR 4/2) fine subangular GRAVEL with medium to coarse Sand |
| 526 | | | | | | GP-SW | | |
| 528 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded medium to coarse SAND |
| 530 | | | | | | CL | | |
| 532 | | | | | | | | Dark gray (10 YR 4/1) coarse Sandy lean CLAY |
| 534 | | | | | | | | |
| 536 | | | | | | | | |
| 538 | | | | | | | | |
| 540 | | 970 | 12 | | | | | |
| 542 | | | | | | | | |
| 544 | | | | | | | | |
| 546 | | | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|---|
| 548 | 30 60 90 | | | | Magothy | SC | | Dark gray (10 YR 4/1) Clayey medium to coarse SAND <i>(continued)</i> |
| 550 | | | | | | SC | | |
| 552 | | | | | | SP | | Light brownish gray (10 YR 6/2) poorly graded medium SAND with trace Silt |
| 554 | | | | | | SP | | |
| 556 | | | | | | SP | | |
| 558 | | | | | | SW | | Light brownish gray (10 YR 6/2) poorly graded medium SAND |
| 560 | | 900 | 10 | | | SW | | |
| 562 | | | | | | SW-SC | | Light gray (10 YR 7/2) well graded fine to medium SAND |
| 564 | | | | | | SW-SC | | |
| 566 | | | | | | SC | | Light gray (10 YR 7/2) well graded fine to coarse SAND with little Clay |
| 568 | | | | | | SC | | |
| 570 | | | | | | SP | | Light gray (10 YR 7/2) well graded fine to coarse SAND with little Clay |
| 572 | | | | | | SP | | |
| 574 | | | | | | SP | | Gray (10 YR 5/1) Clayey well graded fine to medium SAND |
| 576 | | | | | | SP | | |
| 578 | | 1800 | 8.4 | | | SP | | Grayish brown (10 YR 5/2) poorly graded medium SAND |
| 580 | | | | | | SP | | |
| 582 | | | | | | SP | | Light gray (10 YR 7/2) poorly graded medium SAND |
| 584 | | | | | | SP | | |
| 586 | | | | | | SP | | Light gray (10 YR 7/2) poorly graded medium SAND |
| 588 | | | | | | SP | | |
| 590 | | | | | | SP | | Light gray (10 YR 7/2) poorly graded medium SAND |
| 592 | | | | | | SP | | |
| 594 | | | | | | SP | | Light gray (10 YR 7/2) poorly graded medium SAND, few Clay |
| 596 | | | | | | SP | | |
| 598 | | 350 | 0.4 | | | SP | | Very dark gray (10 YR 3/2) well graded fine to medium Clayey SAND |
| 600 | | | | | | SC | | |
| 602 | | | | | | CL | | Very dark gray (10 YR 3/1) lean CLAY with few poorly graded fine Sand |
| 604 | | | | | | | | |
| 606 | | | | | | | | |
| 608 | | | | | | | | |

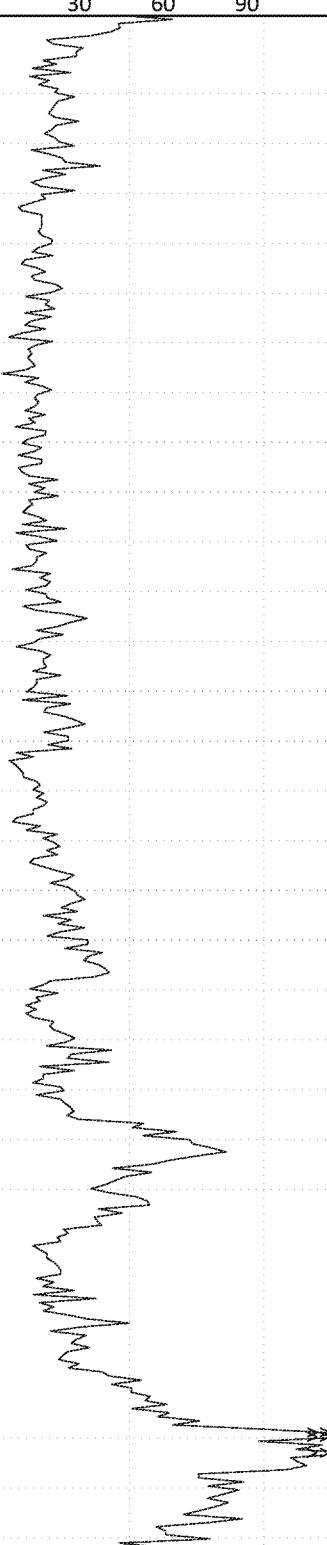
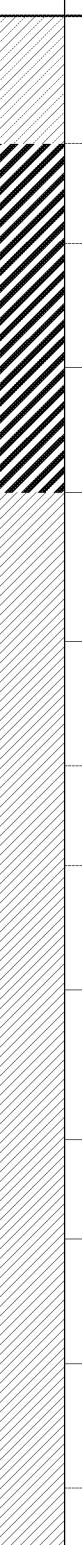
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| DEPTH (ft) | Gamma Ray 30 60 90 | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------------------|-----------|------------|------------|-----------|-------|----------------|--|
| 610 | | | | | Magothy | CL | | Very dark gray (10 YR 3/1) lean CLAY with few poorly graded fine Sand (continued) |
| 612 | | | | | | CH | | Very dark gray (10 YR 3/1) fat CLAY with few poorly graded fine Sand |
| 614 | | | | | | SC | | Dark gray (10 YR 4/1) well graded fine to coarse SAND, with few Clay |
| 616 | | | | | | SC | | Dark Gray (10 YR 4/1) well graded fine to coarse SAND, with little Clay |
| 618 | | | | | | SW | | Dark gray (10 YR 4/1) well graded fine to coarse SAND, few Clay, trace fine gravel |
| 620 | | | | | | SW | | Gray (10 YR 5/1) well graded fine to coarse SAND, few Clay, trace fine gravel |
| 622 | | | | | | SW | | Gray (10 YR 5/1) well graded fine to coarse SAND, few Clay, trace fine gravel |
| 624 | | | | | | SW | | Very pale brown (10 YR 7/3) well graded fine to coarse SAND with interbedded layers of lean Clay (black) |
| 626 | | | | | | SW/CL | | Dark gray (10 YR 4/1) lean CLAY, few poorly graded fine Sand |
| 628 | | | | | | CL | | Gray (10 YR 6/1) well graded fine to coarse SAND, trace lean Clay |
| 630 | | | | | SW | | | Gray (10 YR 6/1) well graded fine to coarse SAND, trace lean Clay |
| 632 | | | | | | | | Gray (10 YR 6/1) well graded fine to coarse SAND, trace lean Clay |
| 634 | | | | | | | | Gray (10 YR 6/1) well graded fine to coarse SAND, trace lean Clay |
| 636 | | | | | | | | Gray (10 YR 6/1) well graded fine to coarse SAND, trace lean Clay |
| 638 | | | | | | | | Light brownish gray (10 YR 6/2) Clayey well graded fine SAND |
| 640 | | | | | 0 | | | |
| 642 | | | | | | | | |
| 644 | | | | | | | | |
| 646 | | | | | 1800 | | | |
| 648 | | | | | | | | |
| 650 | | | | | | | | |
| 652 | | | | | 5.6 | | | |
| 654 | | | | | | | | |
| 656 | | | | | | | | |
| 658 | | | | | 0 | | | |
| 660 | | | | | | | | |
| 662 | | | | | | | | |
| 664 | | | | | 600 | | | |
| 666 | | | | | | | | |
| 668 | | | | | | | | |
| 670 | | | | | < 0.5 | | | |

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| DEPTH (ft) | Gamma Ray 30 60 90 | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------------------------|-----------|------------|------------|-----------|-------|----------------|---|
| 672 | | | | | Magothy | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine SAND (continued) |
| 674 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine SAND |
| 676 | | | | | | SC | | Light brownish gray (10 YR 6/2) Clayey well graded fine SAND |
| 678 | | 550 | 2.8 | | | SW | | Gray (10 YR 6/1) well graded fine to coarse SAND, few lean Clay |
| 680 | | | | | | SW | | Gray (10 YR 6/1) well graded fine to coarse SAND, few lean clay |
| 682 | | | | | | SW | | Light gray (10 YR 7/1) well graded fine to coarse SAND, little lean clay, trace fine gravel |
| 684 | | | | | | CL | | Light gray (10 YR 7/1) well graded fine to coarse Sandy lean CLAY, trace fine subangular Gravel |
| 686 | | | | | | SC | | White (10 YR 8/1) Clayey poorly graded fine SAND, trace fine subangular Gravel |
| 688 | | | | | | SM-SC | | White (10 YR 8/1) Silty Clayey poorly graded fine SAND |
| 690 | | | | | | ML-CL | | White (10 YR 8/1) lean Clayey SILT, few poorly graded fine sand |
| 692 | | | | | | SC | | White (10 YR 8/1) poorly graded coarse Clayey SAND, trace subangular fine Gravel |
| 694 | | | | | | SP/CL | | White (10 YR 8/1) poorly graded coarse SAND with lean Clay |
| 696 | | | | | | CL | | Light brownish gray (10 YR 6/2) lean CLAY with trace well graded fine to coarse Sand |
| 698 | | | | | | | | |
| 700 | | 700 | 4.6 | | | | | |
| 702 | | | | | | | | |
| 704 | | | | | | | | |
| 706 | | | | | | | | |
| 708 | | | | | | | | |
| 710 | | | | | | | | |
| 712 | | | | | | | | |
| 714 | | | | | | | | |
| 716 | | | | | | | | |
| 718 | | | | | | | | |
| 720 | | < 10 | < 10 | | | | | |
| 722 | | | | | | | | |
| 724 | | | | | | | | |
| 726 | | | | | | | | |
| 728 | | | | | | | | |
| 730 | | | | | | | | |
| 732 | | | | | | | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray 30 60 90 | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|--|-----------|------------|------------|-----------|------|--|--|
| 734 |  | | | | Magothy | CL |  | Grayish brown (10 YR 5/2) lean CLAY with few well graded fine to coarse Sand (continued) |
| 736 | | | | | | CL | | |
| 738 | | | | | | CH | | Light gray (10 YR 7/1) fat CLAY with few well graded fine to medium Sand |
| 740 | | 1.4 | < 1.0 | | | CH | | |
| 742 | | | | | | CH | | Light gray (10 YR 7/1) fat CLAY with few well graded fine to medium Sand |
| 744 | | | | | | CH | | |
| 746 | | | | | | CL | | White (10 YR 8/1) fat CLAY with little fine to medium Sand |
| 748 | | | | | | CL | | |
| 750 | | | | | | CL | | White (10 YR 8/1) well graded fine to medium Sandy lean CLAY |
| 752 | | | | | | CL | | |
| 754 | | | | | | CL | | White (10 YR 8/1) poorly graded fine Sandy lean CLAY |
| 756 | | | | | | CL | | |
| 758 | | | | | | CL | | White (10 YR 8/1) lean CLAY with little well graded fine to medium Sand |
| 760 | | | | | | CL | | |
| 762 | | | | | | CL | | White (10 YR 8/1) well graded fine to medium Sandy CLAY |
| 764 | | < 0.5 | < 0.5 | | | CL | | |
| 766 | | | | | | CL | | |
| 768 | | | | | | CL | | White (10 YR 8/1) well graded fine to medium Sandy CLAY |
| 770 | | | | | | CL | | |
| 772 | | | | | | CL | | White (10 YR 8/1) well graded fine to medium Sandy CLAY |
| 774 | | | | | | CL | | |
| 776 | | | | | | CL | | |
| 778 | | | | | | CL | | White (10 YR 8/1) lean CLAY with trace poorly graded fine Sand |
| 780 | | < 1.0 | < 1.0 | | | CL | | |
| 782 | | | | | | CL | | White (10 YR 8/1) lean CLAY with trace poorly graded fine Sand |
| 784 | | | | | | CL | | |
| 786 | | | | | | CL | | White (10 YR 8/1) lean CLAY |
| 788 | | | | | | CL | | |
| 790 | | | | | | CL | | Very dark gray (10 YR 3/1) lean CLAY |
| 792 | | | | | | CL | | |
| 794 | | | | | | CL | | |

(Continued Next Page)

| DEPTH (ft) | Gamma Ray 30 60 90 | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------------------|-----------|------------|------------|-----------|-------|----------------|---|
| 796 | | | | | Magothy | CL | | Very dark gray (10 YR 3/1) lean CLAY (continued) |
| 798 | | | < 0.5 | < 0.5 | | CL | | Gray (10 YR 6/1) lean CLAY with little poorly graded fine Sand |
| 800 | | 0 | | | | SP | | Light gray (10 YR 7/1) poorly graded fine SAND with trace Silt |
| 802 | | | | | | ML | | Light gray (10 YR 7/1) SILT with trace poorly graded fine Sand |
| 804 | | | | | | ML-CL | | Gray (10 YR 6/1) Clayey SILT with few poorly graded fine Sand |
| 806 | | | | | | ML | | Dark gray (10 YR 4/1) SILT with few poorly graded fine Sand and lean clay |
| 808 | | | | | | CL | | Dark gray (10 YR 4/1) lean CLAY with few poorly graded fine Sand and silt |
| 810 | | | | | | CL | | Dark gray (10 YR 4/1) lean CLAY |
| 812 | | | | | | ML-CL | | Dark gray (10 YR 4/1) SILT and lean Clay |
| 814 | | | | | | ML | | Gray (10 YR 5/1) SILT with trace poorly graded fine Sand |
| 816 | | | | | | ML-CL | | Gray (10 YR 6/1) SILT and lean Clay, trace poorly graded fine sand |
| 818 | | | | | | ML-CL | | Gray (10 YR 6/1) SILT and lean Clay, trace poorly graded fine sand |
| 820 | | | | | | ML-CL | | Gray (10 YR 6/1) SILT and lean Clay with little fine to medium well graded sand |

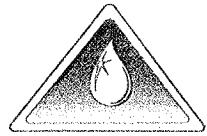
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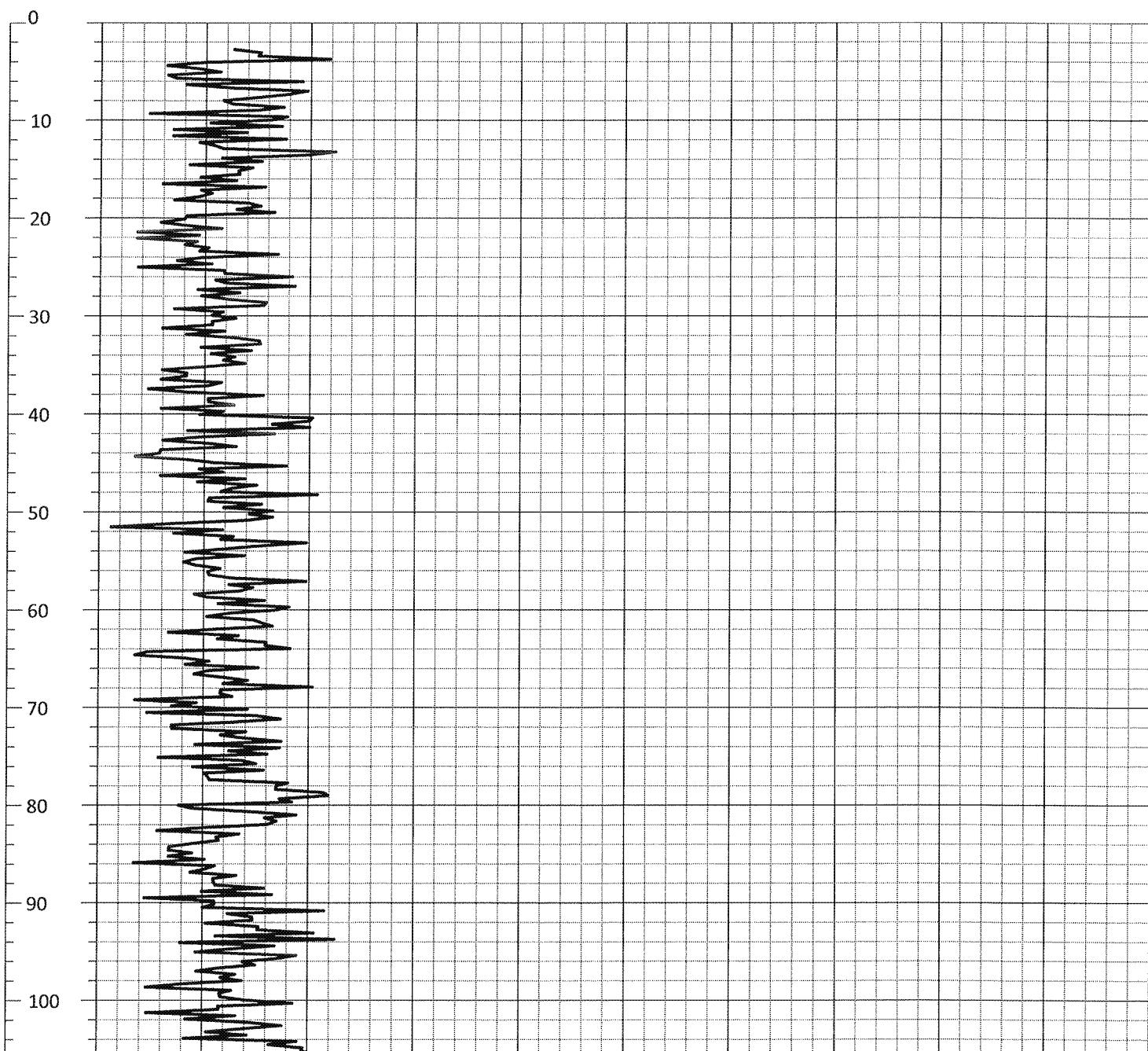
| DEPTH (ft) | Gamma Ray | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | |
|---------------|-----------|-----------|------------|------------|-----------|-------|----------------|---|----|
| | | | | | | | | 30 | 60 |
| 858 | | | | | Magothay | ML-CL | | Gray (10 YR 6/1) SILT with little poorly graded fine Sand and lean clay, trace lignite | |
| 860 | | | | | | | | Gray (10 YR 6/1) SILT with little poorly graded fine Sand and lean clay, trace lignite | |
| 862 | | | | | | | | Gray (10 YR 6/1) SILT with little poorly graded fine Sand and lean clay, trace lignite | |
| 864 | | | | | | | | Gray (10 YR 6/1) Silty well graded fine to coarse SAND with trace fine angular Gravel | |
| 866 | | | | | | | | Gray (10 YR 6/1) Silty well graded fine to coarse SAND with trace fine angular Gravel | |
| 868 | | | | | | | | Very dark gray (10 YR 3/1) well graded medium to coarse SAND with few Silt and lean clay, trace lignite | |
| 870 | | | | | | | | Gray (10 YR 5/1) lean CLAY, trace lignite | |
| 872 | | | | | | | | Gray (10 YR 5/1) lean CLAY, trace lignite | |
| 874 | | | | | | | | Gray (10 YR 5/1) lean CLAY, trace lignite | |
| 876 | | | | | | | | Gray (10 YR 5/1) lean CLAY, trace lignite | |
| 878 | | | | | | | | Gray (10 YR 5/1) lean CLAY, trace lignite | |
| 880 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 882 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 884 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 886 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 888 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 890 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 892 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 894 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 896 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 898 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 900 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 902 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 904 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 906 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 908 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 910 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 912 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 914 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 916 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |
| 918 | | | | | | | | Dark gray (10 YR 4/1) lean CLAY with few lignite | |

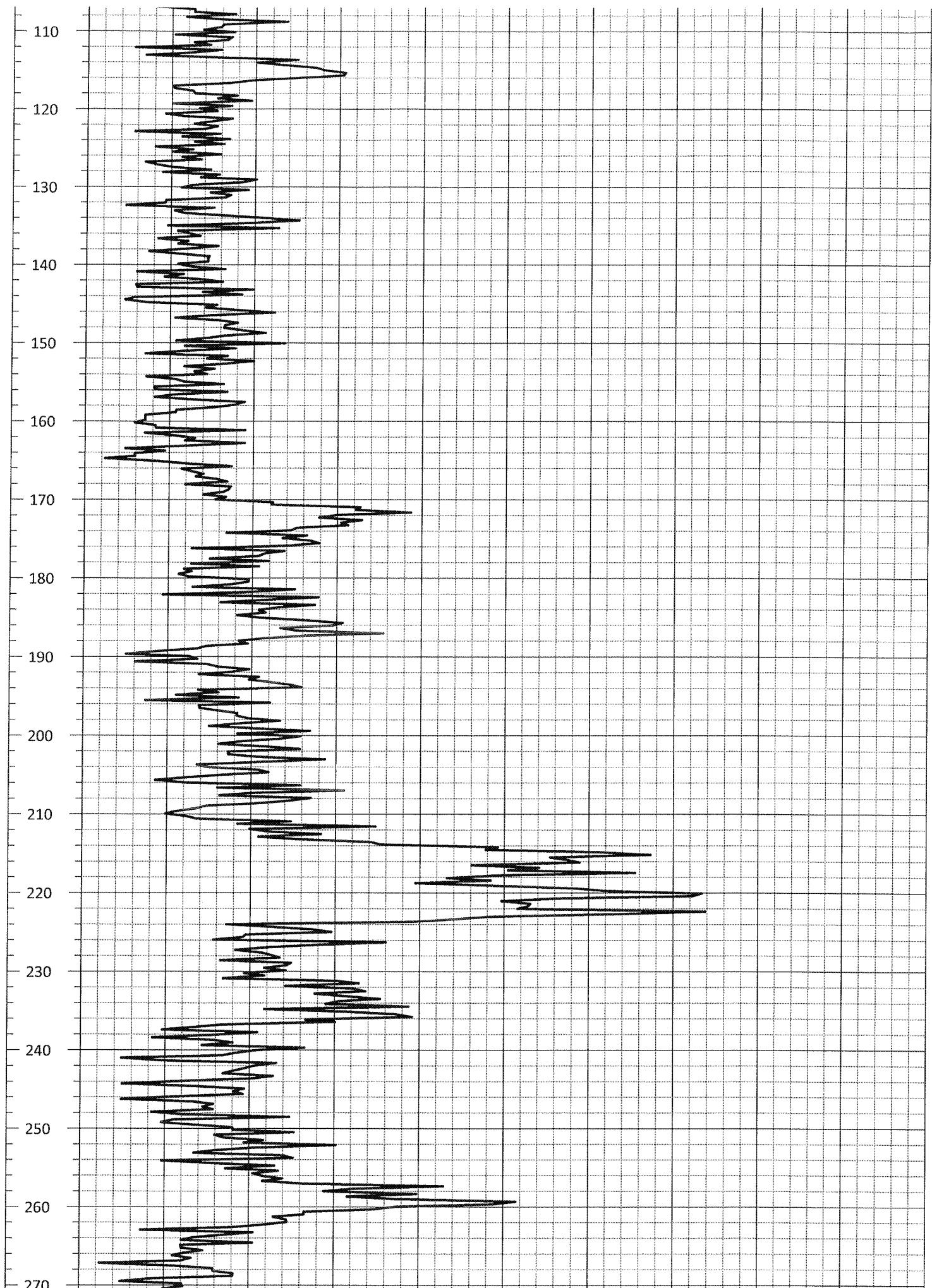
(Continued Next Page)

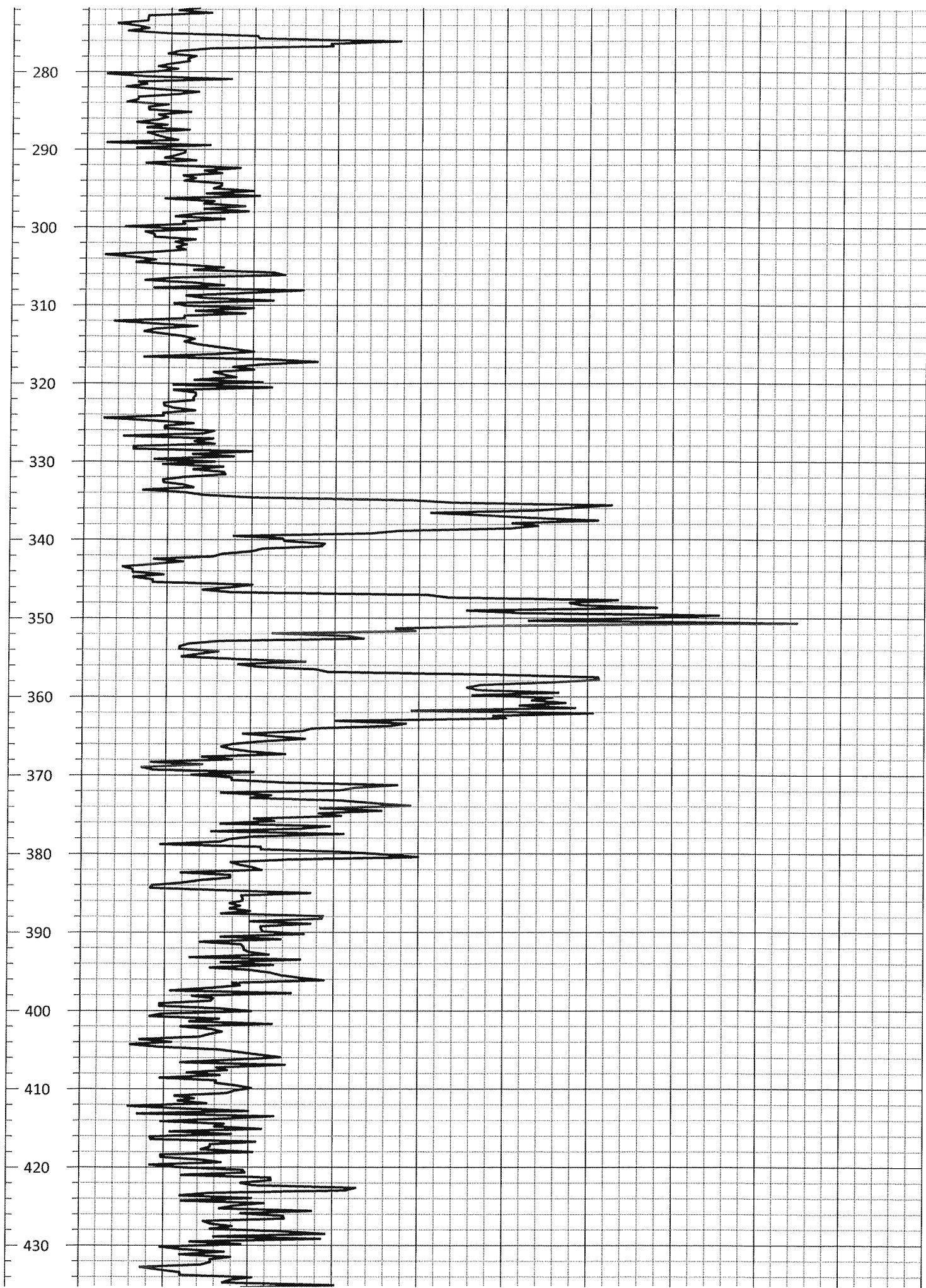
| DEPTH (ft) | Gamma Ray | | | PID (ppm) | TCE (ug/L) | PCE (ug/L) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION |
|---------------|-----------|----|----|-----------|------------|------------|--------------|--------------|----------------|---|
| 918 | 30 | 60 | 90 | | | | | | | |
| 920 | | | | | < 25 | < 25 | Magothy | CL | | Dark gray (10 YR 4/1) lean CLAY with few lignite |
| 922 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 924 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 926 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 928 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 930 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 932 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 934 | | | | | | | | CL | | Mottled red (10 YR 4/6) and gray (10 YR 5/1) lean CLAY |
| 936 | | | | | | | | CH | | Mottled red (10 R 4/6) and brown (10 YR 4/3) fat CLAY with lamination |
| 938 | | | | 0 | | | | Raritan Clay | | Mottled red (10 R 4/6), brown (10 YR 4/3) and gray (10 YR 5/1) fat CLAY with lamination |
| 940 | | | | | | | Raritan Clay | CH | | Mottled red (10 R 4/6) and light gray (10 YR 7/1) fat CLAY with lamination |
| 942 | | | | 0 | | | | CH | | Mottled red (10 R 4/6) and light gray (10 YR 7/1) fat CLAY with lamination |
| 944 | | | | | | | | CH | | Mottled red (10 R 4/6) and light gray (10 YR 7/1) fat CLAY with lamination |
| 946 | | | | | | | | | | |
| 948 | | | | | | | | | | |
| 950 | | | | 0 | | | | | | End of boring at 950.0 ft. bgs. |

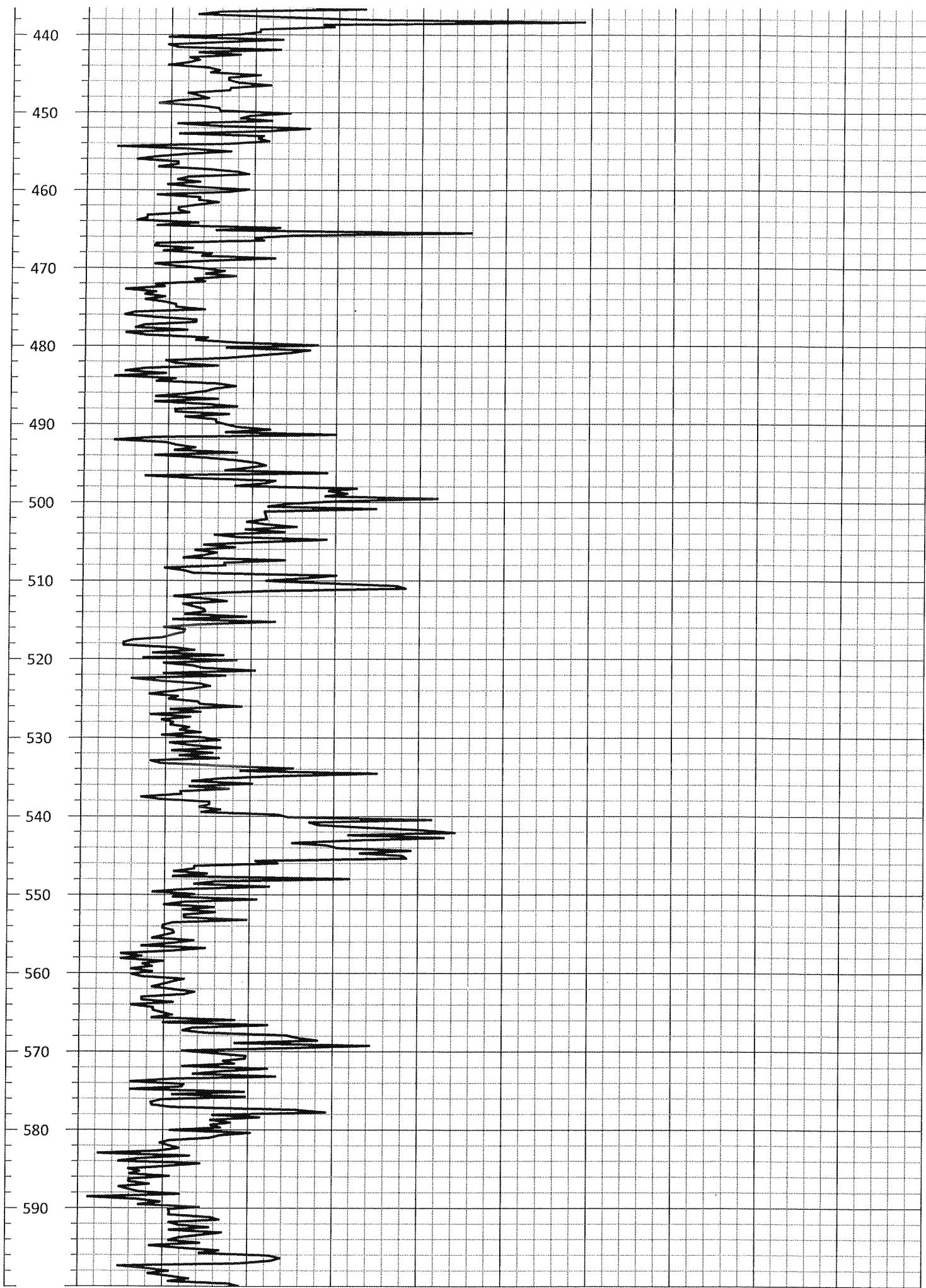
DOWN

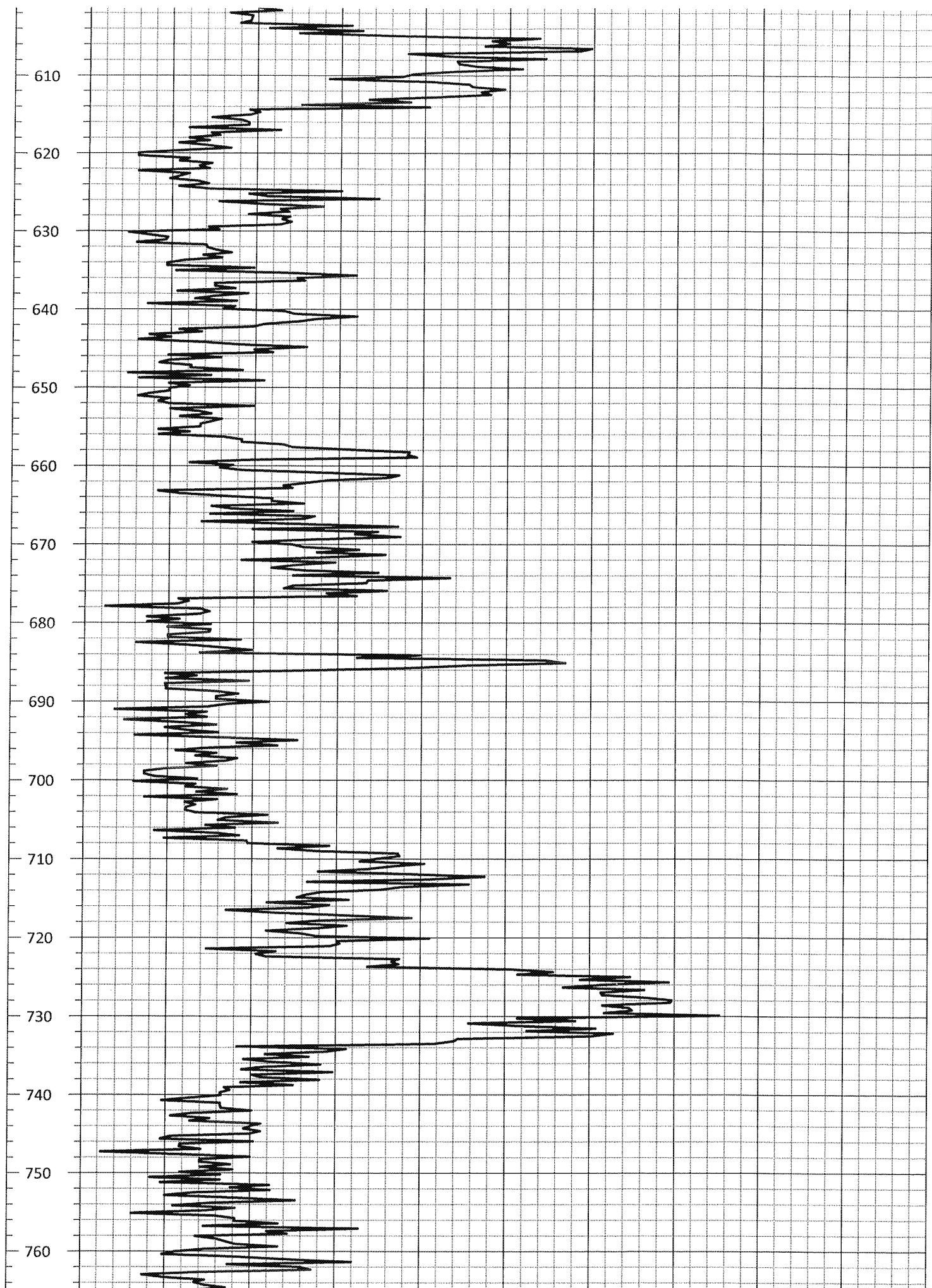
| | | | |
|---|--------------------------------------|----------------|-------|
|  | COMPANY: DELTA WELL & PUMP CO., INC. | | |
| LOCATION: NWIRP SHELLEY DR | | | |
| Well: VP-154 | | Depth Driller: | |
| | | Depth Logger: | |
| Date: 09/04/2014 | Time: | Logged by: CMO | |
| File Name: 739 | | Witness: VIN | |
| Depth (ft.) | <u>0.0</u> | GAMMA (cps) | 100.0 |

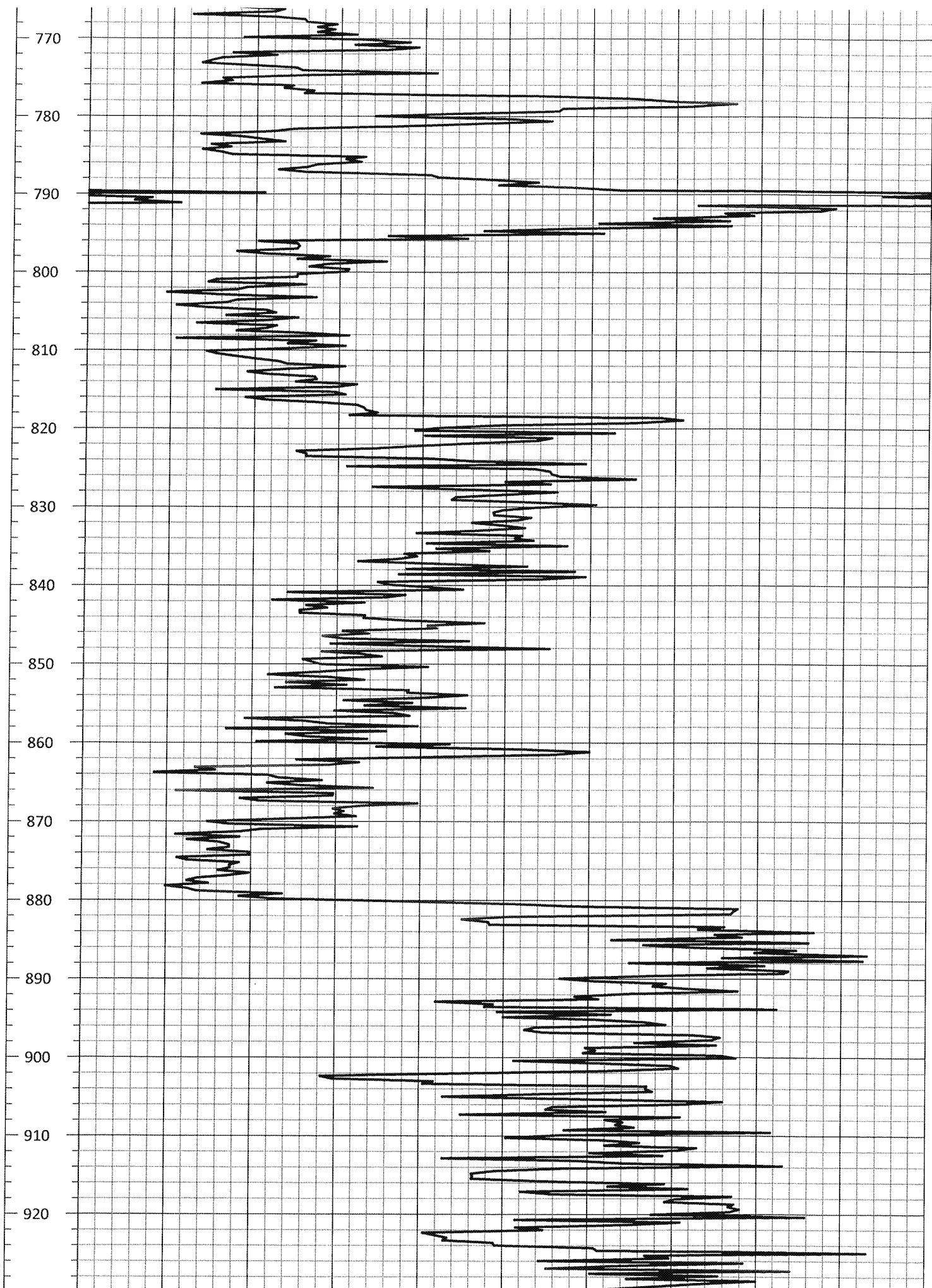


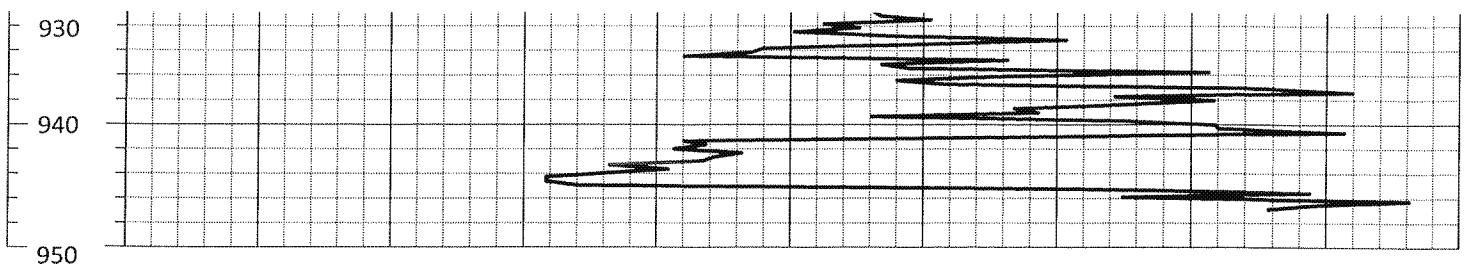








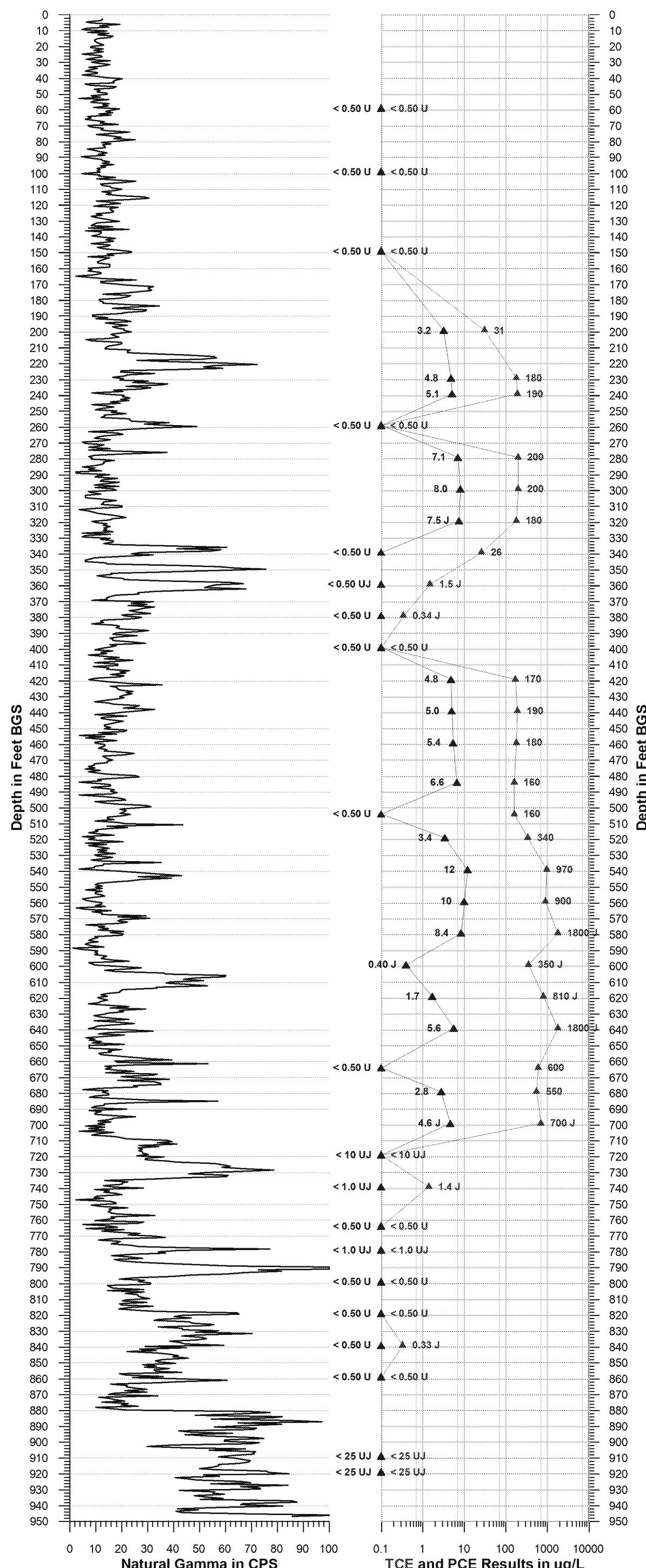




| Depth (ft.) | GAMMA (cps) | |
|-------------|----------------|--|
| 0.0 | 100.0 | |

Section 2
VPB 154 Gamma and PCE/TCE Plot

Vertical Profile Boring VPB-154
Downward Run - September 4, 2014
Validated Analytical Data



Section 3
VPB 154 Groundwater Sample Log Sheets

Client: Navy (ResCon)
Project No: 60266526
Site Location: [REDACTED]
Weather Conds: [REDACTED]

Hydropunch Sample

Navy (ResCon) Date: August 30 1944
 60266526 V.P.B. 154
 (P) - (T) (S) - (F) Dc.
 Collector(s): V.J.

Page 1 of 3

Section 4

VPB 154 Analytical Data Validation

- Analytical Data Sheets
- Chain of Custody Records
- Validation Letter and Table



Resolution Consultants
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fax

Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage

Laboratory: Katahdin Analytical

Service Request: SH6262

Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS)

Validation Level: 3

AECOM Project 60266526.SA.DV

Number:

Prepared by: Dawn Brule/RESCON Completed on: 11/20/2014

Reviewed by: Lori Herberich/RESCON File Name: SH6262_8260B

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 7, 2014 and May 2, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|---|
| VPB154-GW-080714-148-150 | Groundwater |
| VPB154-GWD-080714 | Field Duplicate of VPB154-GW-080714-148-150 |
| VPB154-GW-080714-58-60 | Groundwater |
| VPB154-GW-080714-98-100 | Groundwater |
| VPB154-TB-080714 | Trip Blank |

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW846*, specifically *SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories*, Version 4.2 (DoD, October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✗ Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results

- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID in the report. The submitted EDD file reflects the full sample ID.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recoveries (%Rs) acceptance criteria were met;

- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|--------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R <20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|---|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift >20% | J* | UJ* |
| * No guidance in NFG, thus professional judgment was used | | |

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Table A-3.

Sample results were qualified as follows:

| Blank type | Blank result | Sample result | Action for samples |
|--|----------------------|---------------------------------|--|
| Method, Storage, Field, Trip, or Instrument* | Detects ≤ LOQ | Not detected | No qualification |
| | | < LOQ | Report sample LOQ value with a U |
| | | ≥ LOQ and ≤ 2x LOQ | Report the sample result with a U** |
| | | ≥ 2x the LOQ | No qualifications |
| | > LOQ | < LOQ | Report sample LOQ value with a U |
| | | ≥ LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R |
| | | ≥ LOQ and ≥ blank contamination | If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.** |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment.

LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

Field duplicate RPDs were reviewed for conformance with the QC criterion of ≤30% for aqueous matrices. This criterion applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-080714-148-150 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-080714-58-60 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-080714-98-100 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-GWD-080714 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bt |
| VPB154-GWD-080714 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-TB-080714 | WQ | TRICHLOROFLUOROMETHANE | | 1.0 | UG/L | UJ | c |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|---|--------------------|-----|---------|
| WG148025-7 | 1,1-DICHLOROETHENE | 126 | 80-120% |
| | CARBON DISULFIDE | 129 | 80-120% |
| | ACETONE | 151 | 80-120% |
| | 2-BUTANONE | 125 | 80-120% |
| | 2-HEXANONE | 123 | 80-120% |
| Associated samples: all samples in SDG SH6262 | | | |

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % D | Limits |
|--|-----------------------|-----|-------------|
| WG148068-4 | TRICHLOROFUOROMETHANE | 21 | $\leq 20\%$ |
| Associated sample: VPB154-TB-080714 | | | |
| WG148235-4 | METHYL ACETATE | -25 | $\leq 20\%$ |
| Associated samples: VPB154-GW-080714-58-60, VPB154-GW-080714-98-100, VPB154-GW-080714-148-150, VPB154-GWD-080714 | | | |

Table A-3 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------------|------------------|--------|------|-------|------------------------|
| VPB154-TB-080714 | CARBON DISULFIDE | 0.42 | 0.50 | UG/L | All SDG SH6262 samples |

Attachment B
Qualifier Codes and Explanations

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bt | Trip blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |



600 Technology Way
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Fax: (207) 775-4029

CHAIN of CUSTODY

**PLEASE BEAR DOWN AND
PRINT LEGIBLY IN PEN**

Page _____ of _____

| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |
|--|-------------|--|------------------------------|-------------|--------------------------|
|  | 8/14 1620 | FD EX | | | |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |
| | |  8/14 0950 | | | |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

KAS-COC1

0000009
ORIGINAL

ED 002631A 00004611-00057

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-1RA
Client ID: 154-080714-58-60
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8511.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 7.7 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | J | 0.30 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

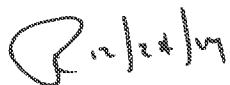
Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6262-1RA
 Client ID: 154-080714-58-60
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6262
 Lab File ID: C8511.D

Sample Date: 07-AUG-14
 Received Date: 09-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.7 | % | | | | | |
| Toluene-d8 | | 95.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 108. | % | | | | | |
| Dibromofluoromethane | | 95.8 | % | | | | | |



R. [Signature]

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-2RA
Client ID: 154-080714-98-100
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8512.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 8.3 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6262-2RA
 Client ID: 154-080714-98-100
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6262
 Lab File ID: C8512.D

Sample Date: 07-AUG-14
 Received Date: 09-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.9 | % | | | | | |
| Toluene-d8 | | 96.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 110. | % | | | | | |
| Dibromofluoromethane | | 97.5 | % | | | | | |

Q12/24/17

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-3RA
Client ID: 154-080714-148-150
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8513.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 12 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.33 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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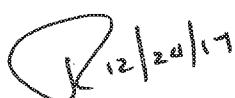
Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6262-3RA
 Client ID: 154-080714-148-150
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6262
 Lab File ID: C8513.D

Sample Date: 07-AUG-14
 Received Date: 09-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.7 | % | | | | | |
| Toluene-d8 | | 96.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 109. | % | | | | | |
| Dibromofluoromethane | | 97.8 | % | | | | | |



J. L. Z. 1/17

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-4RA
Client ID: VPB154-GWD-080714
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8514.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------------------------|--------------------------------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + (S) | 0.28 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 13 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 2.9 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6262-4RA
 Client ID: VPB154-GWD-080714
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6262
 Lab File ID: C8514.D

Sample Date: 07-AUG-14
 Received Date: 09-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | -U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.7 | % | | | | | |
| Toluene-d8 | | 97.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 110. | % | | | | | |
| Dibromofluoromethane | | 100. | % | | | | | |

Q12/24/17

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-5
Client ID: VPB154-TB-080714
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8459.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 12-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148068

Analysis Date: 12-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | J | 0.42 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

12/24/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6262-5
Client ID: VPB154-TB-080714
Project: Navy Clean WE15-03-06 NW
SDG: SH6262
Lab File ID: C8459.D

Sample Date: 07-AUG-14
Received Date: 09-AUG-14
Extract Date: 12-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148068

Analysis Date: 12-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 88.0 | % | | | | | |
| Toluene-d8 | | 90.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 116. | % | | | | | |
| Dibromofluoromethane | | 97.1 | % | | | | | |



Resolution Consultants
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Data Validation Report

| | | |
|-----------------------|---|--------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH6302 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 11/20/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH6302_8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 8 and 11, 2014 and May 2, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-GW-080814-198-200 | Groundwater |
| VPB154-GW-080814-228-230 | Groundwater |
| VPB154-GW-081114-238-240 | Groundwater |
| VPB154-TRIP BLANK-081114 | Trip Blank |

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- ✓ Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicate results

- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Holding Times and Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;

- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|--------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R <20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|---|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift >20% | J* | UJ* |
| * No guidance in NFG, thus professional judgment was used | | |

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required. The QC acceptance criteria were met and/or qualification of the sample results was not required.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

The MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria. 2-Hexanone had high %R for both MS and MSD; however, the associated sample was nondetect for this compound and the results were accepted without qualification.

LCS/LCSD Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|--------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-080814-198-200 | WG | 2-BUTANONE | 20 | 2.5 | UG/L | J | c |
| VPB154-GW-080814-198-200 | WG | ACETONE | 72 | 2.5 | UG/L | J | c |
| VPB154-GW-080814-198-200 | WG | CARBON DISULFIDE | 0.26 | 0.50 | UG/L | J | c |
| VPB154-GW-080814-198-200 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-080814-228-230 | WG | 1,1-DICHLOROETHENE | 1.2 | 0.50 | UG/L | J | c |
| VPB154-GW-080814-228-230 | WG | 2-BUTANONE | 2.7 | 2.5 | UG/L | J | c |
| VPB154-GW-080814-228-230 | WG | ACETONE | 9.8 | 2.5 | UG/L | J | c |
| VPB154-GW-080814-228-230 | WG | CARBON DISULFIDE | 0.25 | 0.50 | UG/L | J | c |
| VPB154-GW-080814-228-230 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-081114-238-240 | WG | 1,1-DICHLOROETHENE | 1.5 | 0.50 | UG/L | J | c |
| VPB154-GW-081114-238-240 | WG | 2-BUTANONE | 3.4 | 2.5 | UG/L | J | c |
| VPB154-GW-081114-238-240 | WG | ACETONE | 16 | 2.5 | UG/L | J | c |
| VPB154-GW-081114-238-240 | WG | CARBON DISULFIDE | 0.30 | 0.50 | UG/L | J | c |
| VPB154-GW-081114-238-240 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | c |
| VPB154-TRIP BLANK-081114 | WQ | METHYL ACETATE | | 0.75 | UG/L | UJ | c |

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|------------|--------------------|-----|---------|
| WG148025-7 | 1,1-DICHLOROETHENE | 126 | 80-120% |
| | CARBON DISULFIDE | 129 | 80-120% |
| | ACETONE | 151 | 80-120% |
| | 2-BUTANONE | 125 | 80-120% |
| | 2-HEXANONE | 123 | 80-120% |

Associated samples: all samples in SDG SH6302

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % R | Limits |
|------------|----------------|-----|--------|
| WG148235-4 | METHYL ACETATE | -25 | <20% |

Associated samples: all samples in SDG SH6302

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |



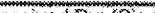
600 Technology Way
Scarborough, ME 04074
Tel: (207) 874-2488
Fax: (207) 775-4029

CHAIN of CUSTODY

**PLEASE BEAR DOWN AND
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Page _____ of _____

COMMENTS

| | | | | | |
|---|-----------------------------|---|------------------------------|-------------|---|
| Relinquished By: (Signature)  | Date / Time 8/14/14 1600 | Received By: (Signature)  | Relinquished By: (Signature) | Date / Time | Received By: (Signature)  8-12-14 0455 |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |
| | | | | | |

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06/1999 9

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6302-1RA
Client ID: VPB154-TB-081114
Project: Navy Clean WE15-03-06 NW
SDG: SH6302
Lab File ID: C8509.D

Sample Date: 11-AUG-14
Received Date: 12-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6302-1RA
Client ID: VPB154-TB-081114
Project: Navy Clean WE15-03-06 NW
SDG: SH6302
Lab File ID: C8509.D

Sample Date: 11-AUG-14
Received Date: 12-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | +/- Q3 | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.5 | % | | | | | |
| Toluene-d8 | | 95.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 106. | % | | | | | |
| Dibromofluoromethane | | 94.0 | % | | | | | |

Giz/za/m

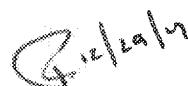
Report of Analytical Results

Client: ENSAFE
Lab ID: SH6302-2RA
Client ID: 154-080814-198-200
Project: Navy Clean WE15-03-06 NW
SDG: SH6302
Lab File ID: C8515.D

Sample Date: 08-AUG-14 **Analysis Date:** 14-AUG-14
Received Date: 12-AUG-14 **Analyst:** DJP
Extract Date: 14-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148235 **Report Date:** 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|--------------------------------|-----------|-------------|-------------|----------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | T | 0.26 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | | 1.7 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 72 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | J | 0.89 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 20 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 31 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 3.2 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Data Validation Report

| | | |
|-----------------------|---|-----------------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH6503 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) and Standard Method 5310 for Total Organic Carbon by High-Temperature Combustion | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/9/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH6503_5310B and 8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 12 and 14, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-FB-081414 | Field blank |
| VPB154-GW-081214-258-260 | Groundwater |
| VPB154-GW-081214-278-280 | Groundwater |
| VPB154-GW-081214-298-300 | Groundwater |
| VPB154-GW-081214-318-320 | Groundwater |
| VPB154-GW-081414-338-340 | Groundwater |
| VPB154-GW-081414-358-360 | Groundwater |
| VPB154-GW-081414-378-380 | Groundwater |
| VPB154-TRIP BLANK-081414 | Trip Blank |

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Standard Methods for the Examination of Water and Wastewater, Method SM5310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and *Quality Systems Manual (QSM)* for Environmental Laboratories, Version 4.2 (DoD, October 2010) where applicable. In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- X Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- X Laboratory blanks/equipment blanks/trip blanks
- X Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- X Laboratory control sample (LCS)/laboratory control sample duplicate (LCSD) results
- NA Field duplicate results
- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (X) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

The 3 vials of sample VPB-154-GW-081414-358-360 each contained mostly soil and not very much liquid. Therefore, each vial was decanted and composited into one vial, then analyzed. Positive and non-detect results for these sample were qualified as estimated (J and UJ) respectively, due to possible loss of sample integrity during the decanting procedure

Holding Times and Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|---------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R < 20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift > 20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method, field, and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Sample results were qualified as follows:

For TOC:

| Blank type | Blank result | Sample result | Action for samples |
|--|---------------|--|---|
| Method, Storage, Field, Trip, or Instrument* | Detects | Not detected | No qualification |
| | $\leq 2x$ LOQ | < 2x LOQ | Report sample LOQ value with a U |
| | | $\geq 2x$ LOQ and \leq 4x LOQ | Report the sample result with a U** |
| | | $>$ 4x LOQ | No qualifications |
| | > 2x LOQ | < 2x LOQ | Report sample LOQ value with a U |
| | | $\geq 2x$ LOQ and < blank contamination | Report the sample result with a U |
| | | $\geq 2x$ LOQ and \geq blank contamination | If the result is $\leq 2x$ blank result, report the sample result U.** If the result is $> 2x$ blank result, no qualification is required.** |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment

[1] Establish an action level (AL) at 5x the blank contamination. If sample result is <AL, qualify the reported result with a U.

For common lab contaminants (methylene chloride, acetone, 2-butanone):

| Blank type | Blank result | Sample result | Action for samples |
|--|---------------|--|--|
| Method, Storage, Field, Trip, or Instrument* | Detects | Not detected | No qualification |
| | $\leq 2x$ LOQ | < 2x LOQ | Report sample LOQ value with a U |
| | | $\geq 2x$ LOQ and \leq 4x LOQ | Report the sample result with a U** |
| | | $>$ 4x LOQ | No qualifications |
| | > 2x LOQ | < 2x LOQ | Report sample LOQ value with a U |
| | | $\geq 2x$ LOQ and < blank contamination | Report the sample result with a U |
| | | $\geq 2x$ LOQ and \geq blank contamination | If the result is $\leq 2x$ blank result, report the sample result U.** |

| Blank type | Blank result | Sample result | Action for samples |
|--|--------------|---------------|---|
| | | | If the result is > 2x blank result, no qualification is required.** |
| * Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L. | | | |
| **Based on professional judgment | | | |

For all other compounds:

| Blank type | Blank result | Sample result | Action for samples | |
|--|---------------------------|---|---|--|
| Method, Storage, Field, Trip, or Instrument* | Detects \leq LOQ | Not detected | No qualification | |
| | | < LOQ | Report sample LOQ value with a U | |
| | | \geq LOQ and \leq 2x LOQ | Report the sample result with a U** | |
| | | \geq 2x the LOQ | No qualifications | |
| | > LOQ | < LOQ | Report sample LOQ value with a U | |
| | | \geq LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R | |
| | | \geq LOQ and \geq blank contamination | If the result is \leq 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.** | |
| * Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L. | | | | |
| **Based on professional judgment. | | | | |

LOQ - Limit of Quantitation. Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Tables A-4 and A-5.

Data qualification on the basis of surrogate recovery nonconformances was as follows:

| Criteria | Action | |
|---------------------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > Upper Limit (UL) | J | No qualification |
| 20% \leq %R \leq Lower Limit (LL) | J | UJ |
| %R < 20% | J | R |

Qualified sample results are shown in Table 1.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS/LCSD Results

The LCS/LCSD %Rs and/or relative percent recoveries (RPDs) were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-6.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|--------------------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) | J | R |
| (LL = lower limit, UL = upper limit) | | |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|----------------------|--------|-------|-------|-----------------------|-------------------|
| VPB154-FB-081414 | WQ | TOTAL ORGANIC CARBON | | 1.0* | MG/L | U | bl |
| VPB154-FB-081414 | WQ | 1,1-DICHLOROETHANE | | 1.0* | UG/L | U | bf |
| VPB154-FB-081414 | WQ | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-FB-081414 | WQ | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-FB-081414 | WQ | ACETONE | | 5.0* | UG/L | UJ | c,bf |
| VPB154-FB-081414 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bf |
| VPB154-FB-081414 | WQ | CHLOROFORM | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-258-260 | WG | 2-BUTANONE | 7.0 | 2.5 | UG/L | J | c |
| VPB154-GW-081214-258-260 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-258-260 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-258-260 | WG | ACETONE | 48 | 2.5 | UG/L | J | l,c |
| VPB154-GW-081214-258-260 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bf |
| VPB154-GW-081214-278-280 | WG | 1,1-DICHLOROETHANE | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-278-280 | WG | 1,1-DICHLOROETHENE | 2.2 | 0.50 | UG/L | J | c |
| VPB154-GW-081214-278-280 | WG | 2-BUTANONE | 3.7 | 2.5 | UG/L | J | c |
| VPB154-GW-081214-278-280 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-278-280 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-278-280 | WG | ACETONE | | 15** | UG/L | UJ | c,bf |
| VPB154-GW-081214-278-280 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bf |
| VPB154-GW-081214-278-280 | WG | CHLOROFORM | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-298-300 | WG | 1,1-DICHLOROETHANE | | 1.2** | UG/L | U | bf |
| VPB154-GW-081214-298-300 | WG | 1,1-DICHLOROETHENE | 3.3 | 0.50 | UG/L | J | c |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-081214-298-300 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-298-300 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-298-300 | WG | ACETONE | | 5.0* | UG/L | UJ | c,bf |
| VPB154-GW-081214-298-300 | WG | CHLOROFORM | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-318-320 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 16 | 0.50 | UG/L | J | s |
| VPB154-GW-081214-318-320 | WG | 1,1-DICHLOROETHANE | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-318-320 | WG | 1,1-DICHLOROETHENE | 2.2 | 0.50 | UG/L | J | s,c |
| VPB154-GW-081214-318-320 | WG | 1,2-DICHLOROETHENE, TOTAL | 2.8 | 1.0 | UG/L | J | s |
| VPB154-GW-081214-318-320 | WG | 2-BUTANONE | 1.9 | 2.5 | UG/L | J | s,c |
| VPB154-GW-081214-318-320 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-318-320 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-081214-318-320 | WG | ACETONE | | 5.0* | UG/L | UJ | s,c,bf |
| VPB154-GW-081214-318-320 | WG | CHLOROFORM | | 1.0* | UG/L | U | bf |
| VPB154-GW-081214-318-320 | WG | CIS-1,2-DICHLOROETHENE | 2.8 | 0.50 | UG/L | J | s |
| VPB154-GW-081214-318-320 | WG | DICHLORODIFLUOROMETHANE | 0.79 | 1.0 | UG/L | J | s |
| VPB154-GW-081214-318-320 | WG | TETRACHLOROETHENE | 7.5 | 0.50 | UG/L | J | s |
| VPB154-GW-081414-338-340 | WG | 1,1-DICHLOROETHANE | | 1.0* | UG/L | U | bf |
| VPB154-GW-081414-338-340 | WG | 1,1-DICHLOROETHENE | 1.5 | 0.50 | UG/L | J | c |
| VPB154-GW-081414-338-340 | WG | 2-BUTANONE | 1.5 | 2.5 | UG/L | J | c |
| VPB154-GW-081414-338-340 | WG | ACETONE | | 5.0* | UG/L | UJ | c,bf |
| VPB154-GW-081414-338-340 | WG | CHLOROFORM | | 1.0* | UG/L | U | bf |
| VPB154-GW-081414-358-360 | WG | 1,1,1-TRICHLOROETHANE | | 0.50 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-081414-358-360 | WG | 1,1,2,2-TETRACHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,1,2-TRICHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,1-DICHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,1 DICHLOROETHENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2,4-TRICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DIBROMOETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DICHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DICHLOROETHENE, TOTAL | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,2-DICHLOROPROPANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,3-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 1,4-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | 2-BUTANONE | 3.8 | 2.5 | UG/L | J | mc,c |
| VPB154-GW-081414-358-360 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | mc,c |
| VPB154-GW-081414-358-360 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | mc,c |
| VPB154-GW-081414-358-360 | WG | ACETONE | | 17** | UG/L | UJ | mc,c,bf |
| VPB154-GW-081414-358-360 | WG | BENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | BROMODICHLOROMETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | BROMOFORM | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-081414-358-360 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | mc,c,bf |
| VPB154-GW-081414-358-360 | WG | CARBON TETRACHLORIDE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CHLOROFORM | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CHLOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CIS-1,2-DICHLOROETHENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CIS-1,3-DICHLOROPROPENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | CYCLOHEXANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | DIBROMOCHLOROMETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | DICHLORODIFLUOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | ETHYLBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | ISOPROPYLBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | M- AND P-XYLENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | METHYL TERT-BUTYL ETHER | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | METHYLENE CHLORIDE | | 2.5 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | O-XYLENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | STYRENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | TETRACHLOROETHENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | TOLUENE | | 0.50 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-081414-358-360 | WG | TRANS-1,2-DICHLOROETHENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | TRANS-1,3-DICHLOROPROPENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | TRICHLOROETHENE | 1.5 | 0.50 | UG/L | J | mc |
| VPB154-GW-081414-358-360 | WG | TRICHLOROFLUOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | VINYL CHLORIDE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-081414-358-360 | WG | XYLENES, TOTAL | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-081414-378-380 | WG | ACETONE | | 5.0* | UG/L | UJ | c,bf |
| VPB154-GW-081414-378-380 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bl |
| VPB154-TRIP BLANK-081414 | WQ | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-081414 | WQ | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-081414 | WQ | ACETONE | | 5.0* | UG/L | UJ | c,bf |
| VPB154-TRIP BLANK-081414 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bf |

*LOQ

**Sample result

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|------------|---|-----|---------|
| WG148025-7 | 1,1-DICHLOROETHENE | 126 | 80-120% |
| | CARBON DISULFIDE | 129 | 80-120% |
| | ACETONE | 151 | 80-120% |
| | 2-BUTANONE | 125 | 80-120% |
| | 2-HEXANONE | 123 | 80-120% |
| | Associated samples: all samples in SDG SH6503 | | |

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % D | Limits |
|--|----------------------|-----|--------|
| WG148353-4 | 4-METHYL-2-PENTANONE | 23 | ≤20% |
| | 2-HEXANONE | 27 | ≤20% |
| Associated samples: VPB154-GW-081214-258-260, VPB154-GW-081214-278-280, VPB154-GW-081214-298-300, VPB154-GW-081214-318-320, VPB154-GW-081414-358-360, VPB154-FB-081414, VPB154-TRIP BLANK-081414 | | | |

Table A-3 - Lab Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------|----------------------|--------|------|-------|--------------------------|
| WG148878-1 | TOTAL ORGANIC CARBON | 0.25 | 0.50 | MG/L | VPB154-FB-081414 |
| WG148424-2 | CARBON DISULFIDE | 0.47 | 0.50 | UG/L | VPB154-GW-081414-378-380 |

Table A-4 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------------|--------------------|--------|------|-------|--|
| VPB154-FB-081414 | 1,1-DICHLOROETHANE | 0.38 | 0.50 | UG/L | VPB154-GW-081214-278-280, VPB154-GW-081214-298-300, VPB154-GW-081214-318-320, VPB154-GW-081414-338-340 |
| VPB154-FB-081414 | ACETONE | 8.2 | 2.5 | UG/L | VPB154-GW-081214-278-280, VPB154-GW-081214-298-300, VPB154-GW-081214-318-320, VPB154-GW-081414-338-340, VPB154-GW-081414-358-360, VPB154-GW-081414-378-380, VPB154-TRIP BLANK-081414 |
| VPB154-FB-081414 | CARBON DISULFIDE | 0.37 | 0.50 | UG/L | VPB154-GW-081214-258-260, VPB154-GW-081214-278-280, VPB154-GW-081414-358-360, VPB154-TRIP BLANK-081414 |
| VPB154-FB-081414 | CHLOROFORM | 1.0 | 0.50 | UG/L | VPB154-GW-081214-278-280, VPB154-GW-081214-298-300, VPB154-GW-081214-318-320, VPB154-GW-081414-338-340 |

Table A-5 - Surrogates

| Sample ID | Surrogate | % Recovery | Lower Limit | Upper Limit |
|--------------------------|-----------------------|------------|-------------|-------------|
| VPB154-GW-081214-318-320 | 1,2-DICHLOROETHANE-D4 | 123 | 70 | 120 |

Table A-6 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|----------|----------------|-------------|-------------|--------------------------|
| WG148353-1 | ACETONE | 153 | 40 | 140 | VPB154-GW-081214-258-260 |

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|--------------------|--|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |
| mc | Method compliance nonconformance |



600 Technology Way
Scarborough, ME 04074
Tel: (207) 874-2400
Fax: (207) 775-4829

CHAIN of CUSTODY

PLEASE BEAR DOWN AND
PRINT LEGIBLY IN PEN

Page _____ of _____

| | | | | | | | | | | | |
|---|-----------------------------|---|---------------------------------|-------------------------------------|--|----------------|----------------|----------------|----------------|----------------|----------------|
| Client Resolution Consultation | | Contact F. Ureche | Phone # (845) 415-4180 | Fax # () | | | | | | | |
| Address 160 Red Schoolhouse Rd | | City Chappaqua | State NY | Zip Code 10511 | | | | | | | |
| Purchase Order # | | Proj. Name / No. NWIRP Nettoze / 66286526 | Katahdin Quote # | | | | | | | | |
| Bill (if different than above) | | Address | | | | | | | | | |
| Sampler (Print / Sign) Vincent Verricchio JV | | Copies To: | | | | | | | | | |
| LAB USE ONLY | WORK ORDER #: | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | | |
| KATAHDIN PROJECT NUMBER SH6503 | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON |
| REMARKS: | | | | | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | | |
| TEMP°C _____ | | <input type="checkbox"/> TEMP BLANK | <input type="checkbox"/> INTACT | <input type="checkbox"/> NOT INTACT | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | | | | | | | |
| | VPD154-6W-081414-318-36 | 8/14/14/94T | 6W | 3 | 3 | | | | | | |
| | VPD154-6W-081414-318-38 | 8/14/14/114T | 6W | 3 | 3 | | | | | | |
| | VPD154-6W-081414-318-39 | 8/14/14/134T | 6W | 3 | 3 | | | | | | |
| | VPD154-6W-081414-318-390 | 8/14/14/1540 | 6W | 3 | 3 | | | | | | |
| | VPB154-6W-081414-318-390 | 8/14/14/1000 | 6W | 3 | 3 | | | | | | |
| | VPB154-6W-081414-318-360 | 8/14/14/1035 | 6W | 3 | 3 | | | | | | |
| | VPD154 - FB - 081414 | 8/14/14/1415 | W | 6 | 3 | 3 | | | | | |
| | VPB154-6W-081414-318-390 | 8/14/14/144T | 6W | 3 | 3 | | | | | | |
| | Trip Bl-K | 5/14/14/1200 | W | 3 | 3 | | | | | | |
| | | / | | | | | | | | | |
| | | / | | | | | | | | | |
| | | / | | | | | | | | | |
| | | / | | | | | | | | | |
| | | / | | | | | | | | | |
| COMMENTS | | | | | | | | | | | |
| Relinquished By: (Signature) <i>Vit Vale</i> | Date / Time 8/14/14 1200 | Received By: (Signature) Fedex | Relinquished By: (Signature) | Date / Time 8/15/14 0915 | Received By: (Signature) <i>Dawn Medeiros</i> | | | | | | |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | | | | | | |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN
SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

000000012

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-1
 Client ID: 154-081214-258-260
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8562.D

Sample Date: 12-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + U J | 0.49 | 1.0 | ug/L | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | - L J | 48 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 7.0 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | + U J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | - L U J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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z/16/15

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-1
Client ID: 154-081214-258-260
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8562.D

Sample Date: 12-AUG-14 **Analysis Date:** 16-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 16-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148353 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.2 | % | | | | | |
| Toluene-d8 | | 97.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 115. | % | | | | | |
| Dibromofluoromethane | | 100. | % | | | | | |

Report of Analytical Results

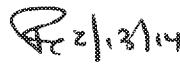
Client: ENSAFE
 Lab ID: SH6503-2
 Client ID: 154-081214-278-280
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8567.D

Sample Date: 12-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|-----------|-------|----------|-----|------------------|----------|---------|
| Dichlorodifluoromethane | J | 0.33 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 2.2 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- U/J | 0.33 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | | 16 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- U/J | 15 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | J | 0.45 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | +/- U | 0.98 1.0 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 4.8 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | +/- U | 0.52 1.0 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | +/- J | 3.7 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | +/- | 2.20 2.00 | ug/L | +/- | 1 | 1.0 4.0 0.28 1.1 | 0.50 2.0 | |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | +/- U/J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 7.1 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | +/- U/J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-2
Client ID: 154-081214-278-280
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8567.D

Sample Date: 12-AUG-14
Received Date: 15-AUG-14
Extract Date: 16-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 4.8 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.0 | % | | | | | |
| Toluene-d8 | | 95.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 120. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-3
 Client ID: 154-081214-298-300
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8568.D

Sample Date: 12-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | J | 0.74 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 3.3 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | | 30 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | + UJ | 4.0 5.0 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 1.2 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 5.2 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | + U | 0.65 1.0 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | E | 230 200 | ug/L | 1 4 | 1 | 1.0 4.0 | 0.28 1.1 | 0.50 2.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | + UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 8.0 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | + UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-3
Client ID: 154-081214-298-300
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8568.D

Sample Date: 12-AUG-14 **Analysis Date:** 16-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 16-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148353 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 5.2 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.1 | % | | | | | |
| Toluene-d8 | | 95.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 115. | % | | | | | |
| Dibromofluoromethane | | 102. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-4
 Client ID: 154-081214-318-320
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8569.D

Sample Date: 12-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | +/- | 0.79 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 2.2 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | J | 16 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- U J | 3.6 5.0 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | +/- U | 0.68 1.0 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | J | 2.8 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | +/- U | 0.81 1.0 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 1.9 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | E | 210 180 | ug/L | +/- | 1 | 10.40 | 0.28 1.1 | 0.56 2.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | +/- U J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | J | 7.5 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | +/- U J | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

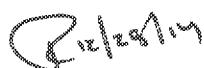
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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-4
Client ID: 154-081214-318-320
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8569.D

Sample Date: 12-AUG-14 **Analysis Date:** 16-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 16-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148353 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | Σ | 2.8 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.0 | % | | | | | |
| Toluene-d8 | | 96.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 123. | % | | | | | |
| Dibromofluoromethane | | 106. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-5RA
Client ID: 154-081414-338-340
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8586.D

Sample Date: 14-AUG-14
Received Date: 15-AUG-14
Extract Date: 18-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148424

Analysis Date: 18-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 1.3 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 1.5 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | | 13 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U J | 87.50 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J U | 0.72 1.0 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.3 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J U | 0.39 1.0 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J U | 1.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | J | 0.34 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 26 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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R.12/29/17

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-5RA
Client ID: 154-081414-338-340
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8586.D

Sample Date: 14-AUG-14 **Analysis Date:** 18-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 18-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148424 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.3 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.5 | % | | | | | |
| Toluene-d8 | | 95.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 117. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-6
 Client ID: 154-081414-358-360
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8559.D

Sample Date: 14-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|-------------|-------------|----------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | | 0.38 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 17 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | | 3.8 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 1.5 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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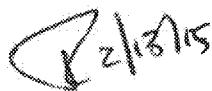
Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-6
 Client ID: 154-081414-358-360
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8559.D

Sample Date: 14-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.3 | % | | | | | |
| Toluene-d8 | | 96.3 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 111. | % | | | | | |
| Dibromofluoromethane | | 98.7 | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-7
 Client ID: VPB154-FB-081414
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8557.D

Sample Date: 14-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + UJ | 0.37 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | + UJ | 8.2 5.0 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | + U | 0.38 1.0 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | + U | 1.0 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | + UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | J | 0.87 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | + UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-7
Client ID: VPB154-FB-081414
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8557.D

Sample Date: 14-AUG-14 **Analysis Date:** 16-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 16-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148353 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|------------------|---------------|--------------|-----------------|------------|----------------|----------------|----------------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.6 | % | | | | | |
| Toluene-d8 | | 97.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 115. | % | | | | | |
| Dibromofluoromethane | | 99.4 | % | | | | | |

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-8RA
Client ID: 154-081414-378-380
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8587.D

Sample Date: 14-AUG-14
Received Date: 15-AUG-14
Extract Date: 18-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148424

Analysis Date: 18-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|-------------|-------------|----------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + | 0.57 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | J | 0.41 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 8.0 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | J | 0.34 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-8RA
Client ID: 154-081414-378-380
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8587.D

Sample Date: 14-AUG-14
Received Date: 15-AUG-14
Extract Date: 18-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148424

Analysis Date: 18-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.5 | % | | | | | |
| Toluene-d8 | | 97.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 112. | % | | | | | |
| Dibromofluoromethane | | 98.2 | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6503-9
 Client ID: VPB154-TB-081414
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6503
 Lab File ID: C8558.D

Sample Date: 14-AUG-14
 Received Date: 15-AUG-14
 Extract Date: 16-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148353

Analysis Date: 16-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|----------------|---------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U S | 0.33* | 1.0 | ug/L | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U S | 4.8-5.0 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U S | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U S | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6503-9
Client ID: VPB154-TB-081414
Project: Navy Clean WE15-03-06 NW
SDG: SH6503
Lab File ID: C8558.D

Sample Date: 14-AUG-14 **Analysis Date:** 16-AUG-14
Received Date: 15-AUG-14 **Analyst:** DJP
Extract Date: 16-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148353 **Report Date:** 19-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.0 | % | | | | | |
| Toluene-d8 | | 95.7 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 110. | % | | | | | |
| Dibromofluoromethane | | 96.7 | % | | | | | |



ANALYTICAL SERVICES

Report of Analytical Results

Client: Rick Purdy
AECOM
701 Edgewater Drive
Wakefield, MA 01880

Lab Sample ID: SH6503-7
Report Date: 29-AUG-14
Client PO: 16518
Project: Navy Clean WE15-03-0
SDG: SH6503

Sample Description
VPB154-FB-081414

| Parameter | Result | Adj LOQ | Adj MDD | Adj LOD | Anal. Method | QC Batch | Anat. Date | Prep. Method | Prep. Date | Footnotes |
|----------------------|------------|---------|---------|---------|--------------|----------|--------------------|--------------|------------|-----------|
| Total Organic Carbon | 0.229 mg/L | 1.0 | 0.10 | .5 | SM5310B | WG148878 | 22-AUG-14 21:59:38 | N/A | N/A | |

1.00

Katahdin Analytical Services 0000237

Cert No E87664

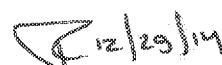


Report of Analytical Results

Client: ENSAFE
Lab ID: SH6302-2RA
Client ID: 154-080814-198-200
Project: Navy Clean WE15-03-06 NW
SDG: SH6302
Lab File ID: C8515.D

Sample Date: 08-AUG-14 **Analysis Date:** 14-AUG-14
Received Date: 12-AUG-14 **Analyst:** DJP
Extract Date: 14-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: DJP **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148235 **Report Date:** 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+p-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | J | 0.89 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 97.0 | % | | | | | |
| Toluene-d8 | | 95.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 110. | % | | | | | |
| Dibromofluoromethane | | 98.5 | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6302-3RA
 Client ID: 154-080814-228-230
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6302
 Lab File ID: C8516.D

Sample Date: 08-AUG-14
 Received Date: 12-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 0.70 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 1.2 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | †† | J | 0.25 | ug/L | 1 | 1 | 0.25 | 0.50 |
| Freon-113 | | 10 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 9.8 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.61 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.4 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | †† J | 2.7 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 180 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 4.8 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | UMM | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6302-3RA
Client ID: 154-080814-228-230
Project: Navy Clean WE15-03-06 NW
SDG: SH6302
Lab File ID: C8516.D

Sample Date: 08-AUG-14
Received Date: 12-AUG-14
Extract Date: 14-AUG-14
Extracted By: DJP
Extraction Method: SW846 5030
Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
Analyst: DJP
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+p-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.6 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.4 | % | | | | | |
| Toluene-d8 | | 95.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 108. | % | | | | | |
| Dibromofluoromethane | | 97.9 | % | | | | | |

G.21/21/14

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6302-4RA
 Client ID: 154-081114-238-240
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6302
 Lab File ID: C8517.D

Sample Date: 11-AUG-14 Analysis Date: 14-AUG-14
 Received Date: 12-AUG-14 Analyst: DJP
 Extract Date: 14-AUG-14 Analysis Method: SW846 8260C
 Extracted By: DJP Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG148235 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|--------------------------------|------------|-------------|-------------|----------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | J | 0.35 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 1.5 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + J | 0.30 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | | 14 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 16 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | | 1.0 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.81 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 4.1 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.86 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.21 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | + J | 3.4 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 190 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 5.1 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

12/29/17

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6302-4RA
 Client ID: 154-081114-238-240
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6302
 Lab File ID: C8517.D

Sample Date: 11-AUG-14
 Received Date: 12-AUG-14
 Extract Date: 14-AUG-14
 Extracted By: DJP
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148235

Analysis Date: 14-AUG-14
 Analyst: DJP
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 16-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 4.1 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 91.4 | % | | | | | |
| Toluene-d8 | | 93.6 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 112. | % | | | | | |
| Dibromofluoromethane | | 98.7 | % | | | | | |

R.2/29/14



Resolution Consultants
250 Apollo Drive
Chelmsford, MA 01824

978.905.2100 tel
978.905.2101 fax

Data Validation Report

| | | |
|-----------------------|---|-----------------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH6621 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) and EPA SW-846 Method 9060A for Total Organic Carbon (TOC) | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/18/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH6621_8260B and 9060A |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 15 and 18, 2014.

| Sample ID | Matrix/Sample Type |
|----------------------------|---|
| VPB154-SOIL-D-081514 | Field Duplicate of VPB154-SOIL-081514-422-424 |
| VPB154-GW-081514-398-400 | Groundwater |
| VPB154-GW-081514-418-420 | Groundwater |
| VPB154-GW-081514-438-440 | Groundwater |
| VPB154-GW-081814-458-460 | Groundwater |
| VPB154-GW-081814-483-485 | Groundwater |
| VPB154-SOIL-081514-422-424 | Soil |
| VPB154-TRIP BLANK-081814 | Trip Blank |

The samples were analyzed in accordance with *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846* (USEPA, 1996), specifically:

- Method 8260B, *Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry*
- Method 9060A, *Total Organic Carbon*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and *Quality Systems Manual (QSM) for Environmental Laboratories*, Version 4.2 (DoD, October 2010) where applicable. In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✗ Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- ✓ Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✗ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Holding Times and Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Table A-1.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift >20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Table A-2.

Sample results were qualified as follows:

| Blank type | Blank result | Sample result | Action for samples |
|--|---------------------------|---|---|
| Method, Storage, Field, Trip, or Instrument* | Detects \leq LOQ | Not detected | No qualification |
| | | < LOQ | Report sample LOQ value with a U |
| | | \geq LOQ and \leq 2x LOQ | Report the sample result with a U** |
| | > LOQ | \geq 2x the LOQ | No qualifications |
| | | < LOQ | Report sample LOQ value with a U |
| | | \geq LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R |
| | | \geq LOQ and \geq blank contamination | If the result is \leq 2x blank result, report the sample result U. If the result is $>$ 2x blank result, no qualification is required.** |
| | | | |
| | | | |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment.

LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

The MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-3.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|--------------------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) | J | R |
| (LL = lower limit, UL = upper limit) | | |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.

Field Duplicate Results

Field duplicate RPDs were reviewed for conformance with the QC criterion of $\leq 60\%$ for soil matrices. This criteria applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-081514-398-400 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-081514-398-400 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-081514-418-420 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 16 | 0.50 | UG/L | J | c,l |
| VPB154-GW-081514-418-420 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-081514-438-440 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 14 | 0.50 | UG/L | J | c,l |
| VPB154-GW-081514-438-440 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-081814-458-460 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 10 | 0.50 | UG/L | J | c,l |
| VPB154-GW-081814-458-460 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-081814-483-485 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 7.7 | 0.50 | UG/L | J | c,l |
| VPB154-TRIP BLANK-081814 | WQ | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 0.50 | UG/L | UJ | c |
| VPB154-TRIP BLANK-081814 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 -Continuing Calibration Verification Standard**

| CCV ID | Compound | % D | Limits |
|---|---------------------------------------|-----|--------|
| WG148554-4 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 31 | <20% |
| Associated samples: all samples in SDG SH6621 | | | |

Table A-2 - Lab Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------|------------------|--------|------|-------|--|
| WG148554-2 | CARBON DISULFIDE | 0.39 | 0.50 | UG/L | VPB154-GW-081514-398-400 VPB154-GW-081514-418-420 VPB154-GW-081514-438-440 VPB154-GW-081814-458-460 VPB154-TRIP BLANK-081814 |

Table A-3 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|---------------------------------------|----------------|-------------|-------------|--|
| WG148554-1 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 140 | 73 | 126 | VPB154-GW-081514-418-420 VPB154-GW-081514-438-440 VPB154-GW-081814-458-460 VPB154-GW-081814-483-485 |

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |



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Page _____ of _____

| | | | | | | | | | | |
|--|---|--|--------------------------|------------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Client <i>Radianz Consultants</i> | Contact <i>E. Vivardon</i> | Phone # <i>(847) 425-4980</i> | Fax # <i>()</i> | | | | | | | |
| Address <i>100 Shel Schoolhouse Rd.</i> | City <i>Christina Ridge</i> | State <i>NY</i> | Zip Code <i>10977</i> | | | | | | | |
| Purchase Order # | Proj. Name / No. <i>NW Reg Dethpage / 66266026</i> | Katahdin Quote # | | | | | | | | |
| Bill (if different than above) | Address | | | | | | | | | |
| Sampler (Print / Sign) <i>Vincent Vercillo</i> | | Copies To: | | | | | | | | |
| LAB USE ONLY | WORK ORDER #: <i>346621</i> | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | |
| KATAHDIN PROJECT NUMBER | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON |
| REMARKS: | | | | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | |
| TEMP'C <input type="checkbox"/> TEMP BLANK <input type="checkbox"/> INTACT <input type="checkbox"/> NOT INTACT | | | | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | <i>VOC</i> | <i>TOL</i> | | | | |
| | <i>UPD154-GW-081514-39866</i> | <i>8/15/14 / 1000</i> | <i>GW</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>UPD154-GW-081514-410-420</i> | <i>8/15/14 / 1145</i> | <i>GW</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>UPD154-Soil-081514-422-424</i> | <i>8/15/14 / 1230</i> | <i>Soil</i> | <i>1</i> | <i>1</i> | | | | | |
| | <i>UPD154-Soil-081514-422-424</i> | <i>8/15/14 / 1230</i> | <i>Soil</i> | <i>1</i> | <i>1</i> | | | | | |
| | <i>UPD154-Soil-D-081514</i> | <i>8/15/14 /</i> | <i>Soil</i> | <i>1</i> | <i>1</i> | | | | | |
| | <i>UPD154-GW-081514-438-440</i> | <i>8/15/14 / 1430</i> | <i>GW</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>UPD154-GW-081814-470-X6</i> | <i>8/18/14 / 1030</i> | <i>GW</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>UPD154-TripBlank-081814</i> | <i>8/18/14 / 1400</i> | <i>W</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>UPD154-GW-081814-482-485</i> | <i>8/18/14 / 1430</i> | <i>GW</i> | <i>3</i> | <i>3</i> | | | | | |
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| COMMENTS | | | | | | | | | | |

| | | | | | |
|---|-------------------------------------|--|------------------------------|-------------|--------------------------|
| Relinquished By: (Signature) <i>Vincent Vercillo</i> | Date / Time <i>8/14/14 16:10</i> | Received By: (Signature) <i>FedEx</i> | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) <i>Clerk 115</i> | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6621-1
 Client ID: 154-081514-398-400
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6621
 Lab File ID: C8637.D

Sample Date: 15-AUG-14
 Received Date: 19-AUG-14
 Extract Date: 20-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|----------------|------------|-------------|----------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | ± 0.034 | 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | UL U/T | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 12 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 2.7 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-1
Client ID: 154-081514-398-400
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8637.D

Sample Date: 15-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.4 | % | | | | | |
| Toluene-d8 | | 95.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 117. | % | | | | | |
| Dibromofluoromethane | | 103. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6621-2
 Client ID: 154-081514-418-420
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6621
 Lab File ID: C8638.D

Sample Date: 15-AUG-14
 Received Date: 19-AUG-14
 Extract Date: 20-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | | 1.9 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | - | 0.33 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | - | 16 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 8.3 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.56 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.87 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.25 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | - | 0.00 | ug/L | - | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 4.8 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-2
Client ID: 154-081514-418-420
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8638.D

Sample Date: 15-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.6 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.8 | % | | | | | |
| Toluene-d8 | | 93.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6621-5
 Client ID: 154-081514-438-440
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6621
 Lab File ID: C8639.D

Sample Date: 15-AUG-14
 Received Date: 19-AUG-14
 Extract Date: 20-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | J | 0.97 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | | 1.7 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- U | 0.34 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | +/- J | 14 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 17 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.49 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 3.0 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.76 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.26 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 4.9 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | E | 240 190 | ug/L | 1 4 | 1 | 1.0 4.0 | 0.28 1.1 | 0.50 2.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 5.0 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-5
Client ID: 154-081514-438-440
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8639.D

Sample Date: 15-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 3.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 95.4 | % | | | | | |
| Toluene-d8 | | 93.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Report of Analytical Results

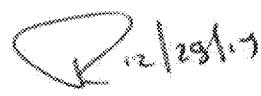
Client: ENSAFE
 Lab ID: SH6621-6
 Client ID: 154-081814-458-460
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6621
 Lab File ID: C8640.D

Sample Date: 18-AUG-14
 Received Date: 19-AUG-14
 Extract Date: 20-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|------------|-------|----------|-----|-----------|------------|------------|
| Dichlorodifluoromethane | J | 0.43 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.28 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | | 1.1 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | ± U | 0.34 ± 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | ± U | 10 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 22 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.44 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.4 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.43 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 4.0 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | ± | 2.0 ± 0.0 | ug/L | ± 4 | 1 | 1.0 ± 0.0 | 0.28 ± 0.1 | 0.50 ± 0.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 5.4 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-6
Client ID: 154-081814-458-460
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8640.D

Sample Date: 18-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|------------|-------|----------|----------|------------|-------------|------------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.4 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 91.6 | % | | | | | |
| Toluene-d8 | | 93.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 103. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6621-7
 Client ID: VPB154-TB-081814
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6621
 Lab File ID: C8631.D

Sample Date: 18-AUG-14
 Received Date: 19-AUG-14
 Extract Date: 20-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|--|--------------------------------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U U U U U | 0.30 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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8/12/2014

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-7
Client ID: VPB154-TB-081814
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8631.D

Sample Date: 18-AUG-14 **Analysis Date:** 20-AUG-14
Received Date: 19-AUG-14 **Analyst:** REC
Extract Date: 20-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148554 **Report Date:** 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.1 | % | | | | | |
| Toluene-d8 | | 96.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 102. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6621-8
Client ID: 154-081814-483-485
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8641.D

Sample Date: 18-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 0.41 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.34 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 0.71 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | L | 7.7 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 9.7 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.24 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 160 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 6.6 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

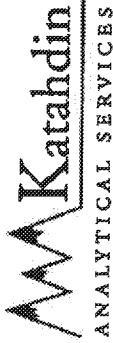
Client: ENSAFE
Lab ID: SH6621-8
Client ID: 154-081814-483-485
Project: Navy Clean WE15-03-06 NW
SDG: SH6621
Lab File ID: C8641.D

Sample Date: 18-AUG-14
Received Date: 19-AUG-14
Extract Date: 20-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148554

Analysis Date: 20-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 21-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | J | 1.6 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.5 | % | | | | | |
| Toluene-d8 | | 94.1 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 116. | % | | | | | |
| Dibromofluoromethane | | 105. | % | | | | | |

ED_002631A_00004611-00148



ANALYTICAL SERVICES

Report of Analytical Results

Client: Rick Purdy
AECOM
701 Edgewater Drive
Wakefield, MA 01880

Lab Sample ID: SH6621-3

Report Date: 05-SEP-14

Client PO: 16518

Project: Navy Clean WE15-03-0

SDG: SH6621

Sample Description

154S-081514-422-424

| Parameter | Result | Adj LOQ | Adj MDL | Adj LOD | Anal. Method | QC Batch | Anal. Date | Prep. Method | Prep. Date | Footnotes |
|--------------|------------------|---------|---------|---------|------------------|----------|--------------------|--------------|------------|-----------|
| TOC In Soil | 1300 ug/g dry wt | \$10 | 110 | 380 | SW846 9060A Mod. | WG148851 | 22-AUG-14 11:55:44 | N/A | N/A | |
| Total Solids | 78. % | 1 | N/A | N/A | SM2540G | WG148631 | 21-AUG-14 10:57:44 | SM2540G | 20-AUG-14 | |

Katahdin Analytical Services 0000188



Cert No E87604



Report of Analytical Results

Client: Rick Purdy
AECOM
701 Edgewater Drive
Wakefield, MA 01880

Lab Sample ID: SH6621-4
Report Date: 05-SEP-14
Client PO: 16518
Project: Navy Clean WE15-03-0
SDG: SH6621

Sample Description
154-SOIL-D-081514

| <u>Sample Description</u> | <u>Matrix</u> | <u>Date Sampled</u> | <u>Date Received</u> |
|---------------------------|---------------|---------------------|----------------------|
| 154-SOIL-D-081514 | SL | 15-AUG-14 | 19-AUG-14 |

| Parameter | Result | Adj LOQ | Adj LOD | Adj MDL | Adj LOD | Anal. Method | QC Batch | Anal. Date | Prep. Method | Prep. Date | Footnotes |
|--------------|----------------|---------|---------|---------|---------------------|--------------|-----------|------------|--------------|------------|-----------|
| TOC In Soil | 2400 ug/gdrywt | 490 | 100 | 370 | SW846 9060A Mod. | WG148851 | 22-AUG-14 | 12:22:24 | N/A | N/A | |
| Total Solids | 82. % | 1. | N/A | N/A | SM7540G | WG148631 | 21-AUG-14 | 10:58:00 | SM2540G | 20-AUG-14 | |

Katahdin Analytical Services 0000189



Cert No E87604



Resolution Consultants
250 Apollo Drive
Chelmsford, MA 01824

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fax

Data Validation Report

| | | |
|-----------------------|---|-----------------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH6777 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) and Standard Method 5310 for Total Organic Carbon by High-Temperature Combustion | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/19/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH6777_5310B and 8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 19 - 21, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|---|
| VPB154-EB-082014 | Equipment blank |
| VPB154-GW-081914-503-505 | Groundwater |
| VPB154-GW-081914-518-520 | Groundwater |
| VPB154-GW-082014-538-540 | Groundwater |
| VPB154-GW-082014-558-560 | Groundwater |
| VPB154-GW-082014-578-580 | Groundwater |
| VPB154-GW-082114-598-600 | Groundwater |
| VPB154-GW-082114-618-620 | Groundwater |
| VPB154-GW-082114-638-640 | Groundwater |
| VPB154-GWD-082114 | Field duplicate of VPB154-GW-082114-638-640 |
| VPB154-TRIP BLANK-082114 | Trip Blank |

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Standard Methods for the Examination of Water and Wastewater, Method SM5310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and Quality Systems Manual (QSM) for Environmental

Laboratories, Version 4.2 (DoD, October 2010) where applicable. In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✗ Laboratory blanks/equipment blanks/trip blanks
- ✗ Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✗ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|---------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R < 20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift > 20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL).

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method, equipment rinsate and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Tables A-3 and A-4.

Sample results were qualified as follows:

For VOC samples:

| Blank type | Blank result | Sample result | Action for samples |
|--|--------------|---|---|
| Method, Storage, Field, Trip, or Instrument* | Detects | Not detected | No qualification |
| | \leq LOQ | < LOQ | Report sample LOQ value with a U |
| | | > LOQ and < 2x LOQ | Report the sample result with a U** |
| | | > 2x the LOQ | No qualifications |
| | > LOQ | < LOQ | Report sample LOQ value with a U |
| | | \geq LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R |
| | | \geq LOQ and \geq blank contamination | If the result is \leq 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.** |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment.

For TOC samples:

| Blank Type | Blank Result | Sample Result | Action for Samples |
|----------------------------|--------------------------|--|---|
| ICB/CCB (Positive) | \geq DL but \leq LOQ | Nondetect | No action |
| | | > DL but < LOQ | Qualify as nondetect (U) at the LOQ |
| | | > LOQ | Use professional judgment (see below [1]) |
| | > LOQ | > DL but < LOQ | Qualify as nondetect (U) at the LOQ |
| | | > LOQ but < ICB/CCB Result | Qualify at level of Blank Result with a "U" or Qualify result as unusable |
| | | > ICB/CCB but < 10x the ICB/CCB result | Qualify as estimated (J) |
| | | > 10x ICB/CCB | No action is taken based on professional judgment |
| PB / EB / FB (Positive) | > LOQ | > DL but < LOQ | Qualify as nondetect (U) at the LOQ |
| | | > LOQ but < 10x Blank Result | Qualify results as unusable |
| | | > 10x Blank Result | No action |
| | \geq DL but \leq LOQ | Nondetect | No action |
| | | > DL but < LOQ | Qualify as nondetect (U) at the LOQ |
| | | > LOQ | Use professional judgment (see below [1]) |

[1] Establish an action level (AL) at 5x the blank contamination. If sample result is < AL, qualify the reported result with a U.
LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-5.

Data qualification on the basis of surrogate recovery nonconformances was as follows:

| Criteria | Action | |
|-----------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > Upper Limit (UL) | J | No qualification |
| 20% ≤ %R < Lower Limit (LL) | J | UJ |
| %R < 20% | J | R |

Qualified sample results are shown in Table 1.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-6.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|---|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) (LL = lower limit, UL = upper limit) | J | R |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.

Field Duplicate Results

Field duplicate RPDs were reviewed for conformance with the QC criterion of $\leq 30\%$ for aqueous matrices. This criterion applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason | |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|-----|
| VPB154-EB-082014 | WQ | TOTAL ORGANIC CARBON | | 1.0* | MG/L | U | bl | |
| VPB154-EB-082014 | WQ | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 0.50 | UG/L | UJ | c | |
| VPB154-EB-082014 | WQ | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c | |
| VPB154-EB-082014 | WQ | BROMOMETHANE | | 1.0 | UG/L | UJ | c | |
| VPB154-EB-082014 | WQ | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c | |
| VPB154-GW-081914-503-505 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 20 | 0.50 | UG/L | J | c,l | |
| VPB154-GW-081914-503-505 | WG | 1,1-DICHLOROETHENE | 2.9 | 0.50 | UG/L | J | c | |
| VPB154-GW-081914-503-505 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c | |
| VPB154-GW-081914-503-505 | WG | ACETONE | 8.4 | 2.5 | UG/L | J | c,l | |
| VPB154-GW-081914-503-505 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c | |
| VPB154-GW-081914-503-505 | WG | CHLOROMETHANE | | 2.0* | UG/L | U | bt | |
| VPB154-GW-081914-503-505 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c | |
| VPB154-GW-081914-518-520 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 26 | 0.50 | UG/L | J | c,l | |
| VPB154-GW-081914-518-520 | WG | 1,1-DICHLOROETHENE | 5.5 | 0.50 | UG/L | J | c | |
| VPB154-GW-081914-518-520 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c | |
| VPB154-GW-081914-518-520 | WG | ACETONE | 8.3 | 2.5 | UG/L | J | c,l | |
| VPB154-GW-081914-518-520 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c | |
| VPB154-GW-081914-518-520 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl | |
| VPB154-GW-081914-518-520 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c | |
| VPB154-GW-082014-538-540 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 61 | 0.50 | UG/L | J | c,l | |
| VPB154-GW-082014-538-540 | WG | 1,1-DICHLOROETHENE | 18 | 0.50 | UG/L | J | c | |
| VPB154-GW-082014-538-540 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c | |
| VPB154-GW-082014-538-540 | WG | 2-BUTANONE | 3.8 | 2.5 | UG/L | J | c | |
| VPB154-GW-082014-538-540 | WG | ACETONE | 20 | 2.5 | UG/L | J | c,l | |
| VPB154-GW-082014-538-540 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c | |
| VPB154-GW-082014-538-540 | WG | CHLOROMETHANE | | 2.0* | UG/L | U | bt | |
| VPB154-GW-082014-538-540 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c | |
| VPB154-GW-082014-558-560 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 57 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082014-558-560 | WG | 1,1-DICHLOROETHENE | | 19 | 0.50 | UG/L | J | c |
| VPB154-GW-082014-558-560 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c | |
| VPB154-GW-082014-558-560 | WG | 2-BUTANONE | 2.4 | 2.5 | UG/L | J | c | |
| VPB154-GW-082014-558-560 | WG | ACETONE | 9.6 | 2.5 | UG/L | J | c,l | |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082014-558-560 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GW-082014-558-560 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082014-578-580 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 120 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082014-578-580 | WG | 1,1-DICHLOROETHENE | 39 | 0.50 | UG/L | J | c |
| VPB154-GW-082014-578-580 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-082014-578-580 | WG | ACETONE | 8.3 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082014-578-580 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GW-082014-578-580 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082014-578-580 | WG | TRICHLOROETHENE | 1800 | 10 | UG/L | J | s |
| VPB154-GW-082114-598-600 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 41 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082114-598-600 | WG | 1,1-DICHLOROETHENE | 14 | 0.50 | UG/L | J | c |
| VPB154-GW-082114-598-600 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-082114-598-600 | WG | 2-BUTANONE | 4.5 | 2.5 | UG/L | J | c |
| VPB154-GW-082114-598-600 | WG | ACETONE | 12 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082114-598-600 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GW-082114-598-600 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-082114-598-600 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082114-598-600 | WG | TRICHLOROETHENE | 350 | 5.0 | UG/L | J | s |
| VPB154-GW-082114-618-620 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 70 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082114-618-620 | WG | 1,1-DICHLOROETHENE | 44 | 0.50 | UG/L | J | c |
| VPB154-GW-082114-618-620 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-082114-618-620 | WG | 2-BUTANONE | 2.3 | 2.5 | UG/L | J | c |
| VPB154-GW-082114-618-620 | WG | ACETONE | 10 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082114-618-620 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GW-082114-618-620 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082114-618-620 | WG | TRICHLOROETHENE | 810 | 5.0 | UG/L | J | s |
| VPB154-GW-082114-638-640 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 69 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082114-638-640 | WG | 1,1-DICHLOROETHENE | 29 | 0.50 | UG/L | J | c |
| VPB154-GW-082114-638-640 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-GW-082114-638-640 | WG | 2-BUTANONE | 1.4 | 2.5 | UG/L | J | c |
| VPB154-GW-082114-638-640 | WG | ACETONE | 6.2 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082114-638-640 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GW-082114-638-640 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082114-638-640 | WG | TRICHLOROETHENE | 1800 | 10 | UG/L | J | s |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GWD-082114 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 70 | 0.50 | UG/L | J | c,l |
| VPB154-GWD-082114 | WG | 1,1-DICHLOROETHENE | 28 | 0.50 | UG/L | J | c |
| VPB154-GWD-082114 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-GWD-082114 | WG | 2-BUTANONE | 1.5 | 2.5 | UG/L | J | c |
| VPB154-GWD-082114 | WG | ACETONE | 6.2 | 2.5 | UG/L | J | c,l |
| VPB154-GWD-082114 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-GWD-082114 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GWD-082114 | WG | TRICHLOROETHENE | 1600 | 10 | UG/L | J | s |
| VPB154-TRIP BLANK-082114 | WQ | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 0.50 | UG/L | UJ | c |
| VPB154-TRIP BLANK-082114 | WQ | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | c |
| VPB154-TRIP BLANK-082114 | WQ | BROMOMETHANE | | 1.0 | UG/L | UJ | c |
| VPB154-TRIP BLANK-082114 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-TRIP BLANK-082114 | WQ | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|------------|--------------------|-----|---------|
| WG148025-7 | 1,1-DICHLOROETHENE | 126 | 80-120% |
| | CARBON DISULFIDE | 129 | 80-120% |
| | ACETONE | 151 | 80-120% |
| | 2-BUTANONE | 125 | 80-120% |
| | 2-HEXANONE | 123 | 80-120% |

Associated samples: all samples in SDG SH6777

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % D | Limits |
|------------|---------------------------------------|-----|-------------|
| WG148777-4 | BROMOMETHANE | -21 | $\leq 20\%$ |
| | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 41 | $\leq 20\%$ |
| | 1,2-DIBROMO-3-CHLOROPROPANE | -23 | $\leq 20\%$ |
| | METHYLCYCLOHEXANE | 26 | $\leq 20\%$ |

Associated samples: all samples in SDG SH6777

Table A-3 - Lab Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------|----------------------|--------|------|-------|--|
| WG148878-1 | TOTAL ORGANIC CARBON | 0.25 | 0.50 | MG/L | VPB154-EB-082014 |
| WG148777-2 | CARBON DISULFIDE | 0.32 | 0.50 | UG/L | VPB154-GW-081914-518-520 VPB154-GW-082114-598-600 VPB154-TRIP BLANK-082114 |

Table A-4 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|--------------------------|---------------|--------|-----|-------|--|
| VPB154-TRIP BLANK-082114 | CHLOROMETHANE | 0.86 | 1.0 | UG/L | VPB154-GW-081914-518-520 VPB154-GW-082114-598-600 |

Table A-5 - Surrogates

| Sample ID | Surrogate | % Recovery | Lower Limit | Upper Limit |
|--------------------------|-----------------------|------------|-------------|-------------|
| VPB154-GW-082014-578-580 | 1,2-DICHLOROETHANE-D4 | 125 | 70 | 120 |
| VPB154-GW-082114-598-600 | 1,2-DICHLOROETHANE-D4 | 126 | 70 | 120 |
| VPB154-GW-082114-618-620 | 1,2-DICHLOROETHANE-D4 | 123 | 70 | 120 |
| VPB154-GW-082114-638-640 | 1,2-DICHLOROETHANE-D4 | 126 | 70 | 120 |
| VPB154-GWD-082114 | 1,2-DICHLOROETHANE-D4 | 126 | 70 | 120 |

Table A-6 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|---------------------------------------|----------------|-------------|-------------|---|
| WG148777-1 | ACETONE | 141 | 40 | 140 | VPB154-GW-081914-503-505 VPB154-GW-081914-518-520 VPB154-GW-082014-538-540 VPB154-GW-082014-558-560 VPB154-GW-082014-578-580 VPB154-GW-082114-598-600 VPB154-GW-082114-618-620 VPB154-GW-082114-638-640 VPB154-GWD-082114 |
| WG148777-1 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 146 | 73 | 126 | VPB154-GW-081914-503-505 VPB154-GW-081914-518-520 VPB154-GW-082014-538-540 VPB154-GW-082014-558-560 VPB154-GW-082014-578-580 VPB154-GW-082114-598-600 VPB154-GW-082114-618-620 VPB154-GW-082114-638-640 VPB154-GWD-082114 |

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bt | Trip blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |



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Page ____ of ____

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|---|---|--|---------------------------------|-------------------------------------|----------------|----------------|----------------|----------------|----------------|--|
| Client <i>Resolution Consultant</i> | Contact <i>E. V. Van der</i> | Phone # <i>(845) 425-4980 ()</i> | Fax # | | | | | | | |
| Address <i>100 Red S. (Hathaway Rd.)</i> | City <i>Clayton Ridge</i> | State <i>NY</i> | Zip Code <i>10977</i> | | | | | | | |
| Purchase Order # | Proj. Name / No. <i>NW11P-Bethpage/166266526</i> | Katahdin Quote # | | | | | | | | |
| Bill (if different than above) | Address | | | | | | | | | |
| Sampler (Print / Sign) <i>Vincent Van Rieloo / U.A.D.</i> | | | Copies To: | | | | | | | |
| LAB USE ONLY | WORK ORDER #: <i>ST16777</i> | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | |
| KATAHDIN PROJECT NUMBER | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | |
| REMARKS: | | <i>Y</i> | <i>T</i> | <i>O</i> | <i>L</i> | <i>Y</i> | <i>O</i> | <i>N</i> | <i>O</i> | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | |
| TEMP°C _____ | | <input type="checkbox"/> TEMP BLANK | <input type="checkbox"/> INTACT | <input type="checkbox"/> NOT INTACT | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | | | | | | |
| | UPR154-(6L-081914-515-105) | 8/19/14 / 1155 | GW | 3 | 3 | | | | | |
| | UPR154-(6L-081914-515-105) | 8/19/14 / 1000 | | | | | | | | |
| | UPR154-GW-081914-518-520 | 8/19/14 / 1460 | GW | 3 | 3 | | | | | |
| | UPR154-GW-081914-538-520 | 8/19/14 / 950 | GW | 3 | 3 | | | | | |
| | UPR154-6L-082014-558-166 | 8/20/14 / 1200 | GW | 3 | 3 | | | | | |
| | UPR154-EB-082014 | 8/20/14 / 1340 | W | 6 | 3 | 3 | | | | |
| | UPR154-GW-082014-578-520 | 8/20/14 / 1415 | GW | 3 | 3 | | | | | |
| | UPR154-6L-082014-598-606 | 8/20/14 / 1000 | GW | 3 | 3 | | | | | |
| | UPR154-GW-082014-618-620 | 8/20/14 / 1210 | GW | 3 | 3 | | | | | |
| | UPR154-GW-082014-618-620 | 8/20/14 / 1415 | GW | 3 | 3 | | | | | |
| | UPR154-Trip blank-082014 | 8/20/14 / 1410 | W | 3 | 3 | | | | | |
| | UPR154-GWD-082014 | 8/20/14 / | GW | 3 | 3 | | | | | |
| | | / | | | | | | | | |
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| COMMENTS | | | | | | | | | | |

| | | | | | |
|--|-----------------------------|--|------------------------------|-------------|---|
| Relinquished By: (Signature) <i>V. Van Rieloo</i> | Date / Time 8/20/14 1620 | Received By: (Signature) <i>Fedex</i> | Relinquished By: (Signature) | Date / Time | Received By: (Signature) <i>8/22/14 0900</i> |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN
SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-5
 Client ID: VPB154-EB-082014
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8691.D

Sample Date: 20-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Report of Analytical Results

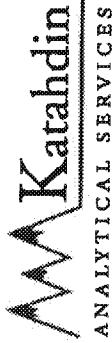
Client: ENSAFE
 Lab ID: SH6777-5
 Client ID: VPB154-EB-082014
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8691.D

Sample Date: 20-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.7 | % | | | | | |
| Toluene-d8 | | 95.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 112. | % | | | | | |
| Dibromofluoromethane | | 102. | % | | | | | |





ANALYTICAL SERVICES

Report of Analytical Results

Client: Rick Purdy
AECOM
701 Edgewater Drive
Wakefield, MA 01880

Lab Sample ID: SH6777-5

Report Date: 09-SEP-14

Client PO: 16518

Project: Navy Clean WE15-03-0

SDG: SH6777

Sample Description

VPB154-EB-082204

Matrix

AQ

Date Sampled

20-AUG-14

Date Received

22-AUG-14

| Parameter | Result | Adj LOQ | Adj MDL | Adj LOD | Anal Method | QC Batch | Anal Date | Prep. Method | Prep. Date | Footnotes |
|----------------------|------------|---------|---------|---------|-------------|----------|--------------------|--------------|------------|-----------|
| Total Organic Carbon | 10.23 mg/l | 1.0 | 0.10 | .5 | SM5310B | WG148878 | 22-AUG-14 22:49:41 | N/A | N/A | |

1.0
0.10
.5

SM5310B
WG148878

22-AUG-14 22:49:41
N/A
N/A

✓ ✓ ✓

✓ ✓ ✓

Katahdin Analytical Services 0000328



Cert No E87604

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-1
Client ID: 154-081914-503-505
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8692.D

Sample Date: 19-AUG-14
Received Date: 22-AUG-14
Extract Date: 22-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | | 2.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | + U | 0.53 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | + U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 2.9 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | + J | 20 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | + J | 8.4 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.2 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.4 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.53 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | | 160 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | J | 0.40 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

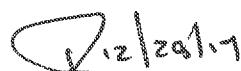
07/28/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-1
Client ID: 154-081914-503-505
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8692.D

Sample Date: 19-AUG-14 **Analysis Date:** 22-AUG-14
Received Date: 22-AUG-14 **Analyst:** REC
Extract Date: 22-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148777 **Report Date:** 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | +/- Q3 | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.2 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | +/- Q3 | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.9 | % | | | | | |
| Toluene-d8 | | 94.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 115. | % | | | | | |
| Dibromofluoromethane | | 101. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-2
 Client ID: 154-081914-518-520
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8693.D

Sample Date: 19-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|---------|----------|
| Dichlorodifluoromethane | J | 0.66 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | +/- U/J | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 5.5 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- U | 0.35 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | +/- J | 26 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- J | 8.3 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.79 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.3 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.74 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.41 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | +/- | 350 340 | ug/L | +/- | 1.0 | 1.0 | 0.2822 | 0.50 4.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | J | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 3.4 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-2
Client ID: 154-081914-518-520
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8693.D

Sample Date: 19-AUG-14 **Analysis Date:** 22-AUG-14
Received Date: 22-AUG-14 **Analyst:** REC
Extract Date: 22-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148777 **Report Date:** 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U or U/T | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.3 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U or U/T | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.5 | % | | | | | |
| Toluene-d8 | | 95.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 113. | % | | | | | |
| Dibromofluoromethane | | 100. | % | | | | | |



R. R. / 2014

Report of Analytical Results

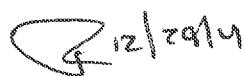
Client: ENSAFE
 Lab ID: SH6777-3
 Client ID: 154-082014-538-540
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8694.D

Sample Date: 20-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|------------|-------|----------|-----|----------|----------|-----------|
| Dichlorodifluoromethane | J | 0.49 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | +/- U | 0.54 ± 0.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | +/- U S | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.42 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 18 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | +/- J | 61 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- J | 20 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 2.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 4.4 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.1 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.87 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | +/- J | 3.8 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | X | 860 ± 70 | ug/L | X 10 | 1 | 1.0 ± 10 | 0.28 ± 8 | 0.50 ± 50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.1 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 12 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2



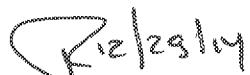
Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-3
Client ID: 154-082014-538-540
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8694.D

Sample Date: 20-AUG-14
Received Date: 22-AUG-14
Extract Date: 22-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 4.4 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.6 | % | | | | | |
| Toluene-d8 | | 93.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 115. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |



Riz/29/14

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-4
 Client ID: 154-082014-558-560
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8695.D

Sample Date: 20-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|---------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | J | 0.49 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | + WS | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.47 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | - J | 19 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | + - J | 57 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | + - J | 9.6 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 2.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 3.9 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.98 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 1.1 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | + - J | 2.4 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | + - | 8.6 9.0 | ug/L | + 10 | 1 | 4.0 10 | 0.28 2.8 | 0.50 5.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | J | 0.98 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 10 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | | 6.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |



Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-4
Client ID: 154-082014-558-560
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8695.D

Sample Date: 20-AUG-14 **Analysis Date:** 22-AUG-14
Received Date: 22-AUG-14 **Analyst:** REC
Extract Date: 22-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148777 **Report Date:** 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 3.9 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.7 | % | | | | | |
| Toluene-d8 | | 94.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 117. | % | | | | | |
| Dibromofluoromethane | | 105. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-6
 Client ID: 154-082014-578-580
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8696.D

Sample Date: 20-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|-----------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 0.79 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | ← U ← J | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.80 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | ← J | 39 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | ← J | 120 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | ← J | 8.3 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 3.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 5.2 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.2 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 2.6 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | ← J | 1200 1800 | ug/L | 1/20 | 1 | 10.20 | 0.2856 | 0.50 10 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.8 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 8.4 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-6
Client ID: 154-082014-578-580
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8696.D

Sample Date: 20-AUG-14
Received Date: 22-AUG-14
Extract Date: 22-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | ±± ug/l | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 5.2 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | ±± ug/l | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 91.7 | % | | | | | |
| Toluene-d8 | | 93.6 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 116. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Riz Lazura

Report of Analytical Results

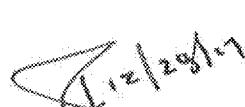
Client: ENSAFE
 Lab ID: SH6777-7
 Client ID: 154-082114-598-600
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8697.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------|--------------|-------------|-------------|-------------|----------|------------|-------------|-------------|
| Dichlorodifluoromethane | J | 0.87 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | + U T | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | T | 14 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + U | 0.27 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | + U T | 41 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | + T | 12 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 1.8 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.2 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 1.3 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | + T | 4.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | + T | 300 | ug/L | X 10 | 1 | 10 | 0.28 | 2.8 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.1 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | J | 0.40 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-7
 Client ID: 154-082114-598-600
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8697.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | UL | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.2 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | UL | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 96.8 | % | | | | | |
| Toluene-d8 | | 97.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 120. | % | | | | | |
| Dibromofluoromethane | | 108. | % | | | | | |

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Report of Analytical Results

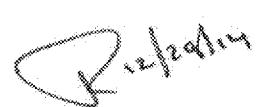
Client: ENSAFE
 Lab ID: SH6777-8
 Client ID: 154-082114-618-620
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8698.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 0.54 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.37 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 44 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | J | 70 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 10 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 4.7 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 3.1 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.5 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 3.7 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 2.3 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | J | 200 | ug/L | 10 | 1 | 10 | 0.28 | 2.5 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 1.7 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-8
Client ID: 154-082114-618-620
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8698.D

Sample Date: 21-AUG-14 **Analysis Date:** 22-AUG-14
Received Date: 22-AUG-14 **Analyst:** REC
Extract Date: 22-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148777 **Report Date:** 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 3.1 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.4 | % | | | | | |
| Toluene-d8 | | 96.1 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 106. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-9
 Client ID: 154-082114-638-640
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8699.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|---------------------------|-------|----------|-----|------------------|---------------------|--------------------|
| Dichlorodifluoromethane | J | 0.39 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | ✓ U ✓ | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.40 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | ✓ J | 29 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | ✓ J | 69 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | ✓ J | 6.2 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 4.5 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 4.4 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.3 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 2.5 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | ✓ J | 1.4 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | ✓ J | 1200 ¹⁸⁰⁰ ug/L | | X 10 | 1 | 10 ²⁰ | 0.28 ^{5.4} | 0.50 ¹⁰ |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 5.6 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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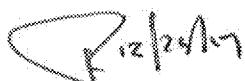
Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-9
Client ID: 154-082114-638-640
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8699.D

Sample Date: 21-AUG-14
Received Date: 22-AUG-14
Extract Date: 22-AUG-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+p-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 4.4 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 91.8 | % | | | | | |
| Toluene-d8 | | 92.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 116. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-11
 Client ID: VPB154-GWD-082114
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8700.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | J | 0.34 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | J | 0.42 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 28 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | J | 70 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 6.2 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | | 4.4 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 4.1 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.3 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | | 2.5 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J | 1.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | J | 1200 | ug/L | 1 | 10 | 10.0 | 0.2856 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 1.6 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 5.6 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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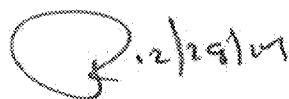
12/25/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6777-11
Client ID: VPB154-GWD-082114
Project: Navy Clean WE15-03-06 NW
SDG: SH6777
Lab File ID: C8700.D

Sample Date: 21-AUG-14 **Analysis Date:** 22-AUG-14
Received Date: 22-AUG-14 **Analyst:** REC
Extract Date: 22-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG148777 **Report Date:** 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 4.1 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 92.1 | % | | | | | |
| Toluene-d8 | | 94.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 107. | % | | | | | |



Q.2/28/14

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-10
 Client ID: VPB154-TB-082114
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8690.D

Sample Date: 21-AUG-14
 Received Date: 22-AUG-14
 Extract Date: 22-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG148777

Analysis Date: 22-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | -+ Q | 0.86 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U+ UJ | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | J | 0.25 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U+ QJ | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U+ W | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6777-10
 Client ID: VPB154-TB-082114
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6777
 Lab File ID: C8690.D

Sample Date: 21-AUG-14 Analysis Date: 22-AUG-14
 Received Date: 22-AUG-14 Analyst: REC
 Extract Date: 22-AUG-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG148777 Report Date: 26-AUG-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+p-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.8 | % | | | | | |
| Toluene-d8 | | 96.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 112. | % | | | | | |
| Dibromofluoromethane | | 103. | % | | | | | |



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Data Validation Report

| | | |
|-----------------------|---|--------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH6880 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/05/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH6880_8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 22 and 25, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-GW-082214-663-665 | Groundwater |
| VPB154-GW-082214-678-680 | Groundwater |
| VPB154-GW-082514-698-700 | Groundwater |
| VPB154-GW-082514-718-720 | Groundwater |
| VPB154-GW-082514-738-740 | Groundwater |
| VPB154-TRIP BLANK-082514 | Trip Blank |

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- Data completeness (chain-of-custody [COC])/sample integrity
- Holding times and sample preservation
- GC/MS performance checks
- Initial calibration/continuing calibration verification
- Laboratory blanks/trip blanks/equipment blanks
- Surrogate spike recoveries
- Matrix spike (MS) and/or matrix spike duplicate (MSD) results

- X Laboratory control sample (LCS) results
- NA Field duplicate results
- ✓ Internal standards results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (X) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Selected samples were mostly soil and had very little standing water.

- For sample VPB154-GW-082514-698-700 the laboratory decanted the liquid from one vial prior to analysis.
- For both samples VPB154-GW-082514-718-720 and VPB154-GW-082514-738-740, the laboratory decanted the water from three individual vials into one vial as a composite. Due to limited volume, the samples were analyzed at dilutions of 1:20 and 1:2, respectively.

Positive and nondetect results for these sample were qualified as estimated (J and UJ, respectively) due to possible loss of sample integrity during the decanting procedure. Qualified sample results are shown in Table 1.

Holding Times and Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|--------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R <20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift >20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Table A-3.

Sample results were qualified as follows:

| Blank type | Blank result | Sample result | Action for samples |
|--|--------------|---------------------------------|--|
| Method, Storage, Field, Trip, or Instrument* | Detects | Not detected | No qualification |
| | | < LOQ | Report sample LOQ value with a U |
| | | ≥ LOQ and ≤ 2x LOQ | Report the sample result with a U** |
| | > LOQ | > 2x the LOQ | No qualifications |
| | | < LOQ | Report sample LOQ value with a U |
| | | ≥ LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R |
| | | ≥ LOQ and ≥ blank contamination | If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.** |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment.

LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-4.

Data qualification on the basis of surrogate recovery nonconformances was as follows:

| Nonconformance | Action | |
|-----------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > Upper Limit (UL) | J | No qualification |
| 20% ≤ %R < Lower Limit (LL) | J | UJ |
| %R < 20% | J | R |

Qualified sample results are shown in Table 1.

MS/MSD Results

The MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-5.

Data qualification to the analytes associated with the specific MS/MSD nonconformances was as follows:

| Nonconformance | Action | |
|------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > UL | J | No qualification |
| 20% ≤ %R < LL | J | UJ |
| %R < 20% (see note 1) | J | R* |
| %RPD > UL (see note 2) | J | No qualification |

Note: Actions are applied to the native unspiked sample only (see note 3)
 *When the native sample concentration is >4X the concentration of the spike added (based on Region I criteria), evaluate the MS, MSD, and native sample with regards to %RSD rather than %R (professional judgment)

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject (R) sample results previously negated (U) on the basis of blank contamination.
2. In the absence of Region 2 guidance, RPD actions are based on professional judgment.
3. If a field duplicate sample was also collected for the native sample chosen for MS/MSD analysis, professional judgment is used to apply MS/MSD actions to the corresponding field duplicate sample as well as the native sample.

Qualified sample results are shown in Table 1.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-6.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|--------------------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) | J | R |
| (LL = lower limit, UL = upper limit) | | |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082214-663-665 | WG | 1,1,1-TRICHLOROETHANE | 2.0 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 84 | 0.50 | UG/L | J | c,s,l |
| VPB154-GW-082214-663-665 | WG | 1,1,2-TRICHLOROETHANE | 1.2 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | 1,1-DICHLOROETHANE | 2.6 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | 1,1-DICHLOROETHENE | 16 | 0.50 | UG/L | J | c,s |
| VPB154-GW-082214-663-665 | WG | 1,2-DICHLOROETHENE, TOTAL | 3.9 | 1.0 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | 2-BUTANONE | 2.4 | 2.5 | UG/L | J | c,s |
| VPB154-GW-082214-663-665 | WG | ACETONE | 10 | 2.5 | UG/L | J | c,s |
| VPB154-GW-082214-663-665 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bt |
| VPB154-GW-082214-663-665 | WG | CARBON TETRACHLORIDE | 2.4 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | CHLOROFORM | 1.6 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | CIS-1,2-DICHLOROETHENE | 3.9 | 0.50 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | DICHLORODIFLUOROMETHANE | 0.26 | 1.0 | UG/L | J | s |
| VPB154-GW-082214-663-665 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082214-678-680 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 21 | 0.50 | UG/L | J | c,l |
| VPB154-GW-082214-678-680 | WG | 1,1-DICHLOROETHENE | 3.2 | 0.50 | UG/L | J | c |
| VPB154-GW-082214-678-680 | WG | ACETONE | 8.7 | 2.5 | UG/L | J | c |
| VPB154-GW-082214-678-680 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | c,bt |
| VPB154-GW-082214-678-680 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | c |
| VPB154-GW-082514-698-700 | WG | 1,1,1-TRICHLOROETHANE | 0.60 | 0.50 | UG/L | J | mc,s,m |
| VPB154-GW-082514-698-700 | WG | 1,1,2,2-TETRACHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 30 | 0.50 | UG/L | J | mc,c,s,l,m |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-----------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082514-698-700 | WG | 1,1,2-TRICHLOROETHANE | 0.68 | 0.50 | UG/L | J | mc,s |
| VPB154-GW-082514-698-700 | WG | 1,1-DICHLOROETHANE | 1.5 | 0.50 | UG/L | J | mc,s,m |
| VPB154-GW-082514-698-700 | WG | 1,1-DICHLOROETHENE | 6.0 | 0.50 | UG/L | J | mc,c,s,m |
| VPB154-GW-082514-698-700 | WG | 1,2,4-TRICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 0.75 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,2-DIBROMOETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,2-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,2-DICHLOROETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,2-DICHLOROETHENE, TOTAL | 2.9 | 1.0 | UG/L | J | mc,s,m |
| VPB154-GW-082514-698-700 | WG | 1,2-DICHLOROPROPANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,3-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 1,4-DICHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 2-BUTANONE | 5.3 | 2.5 | UG/L | J | mc,c,s |
| VPB154-GW-082514-698-700 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | 4-METHYL-2-PENTANONE | | 2.5 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | ACETONE | 26 | 2.5 | UG/L | J | mc,c,s |
| VPB154-GW-082514-698-700 | WG | BENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | BROMODICHLOROMETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | BROMOFORM | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | BROMOMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | CARBON DISULFIDE | | 1.0* | UG/L | UJ | mc,c,bt |
| VPB154-GW-082514-698-700 | WG | CARBON TETRACHLORIDE | | 0.50 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082514-698-700 | WG | CHLOROBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | CHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | CHLOROFORM | 1.2 | 0.50 | UG/L | J | mc,s |
| VPB154-GW-082514-698-700 | WG | CHLOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | CIS-1,2-DICHLOROETHENE | 2.9 | 0.50 | UG/L | J | mc,s,m |
| VPB154-GW-082514-698-700 | WG | CIS-1,3-DICHLOROPROPENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | CYCLOHEXANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | DIBROMOCHLOROMETHANE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | DICHLORODIFLUOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | ETHYLBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | ISOPROPYLBENZENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | M- AND P-XYLENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | METHYL ACETATE | | 0.75 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | METHYL CYCLOHEXANE | | 0.50 | UG/L | UJ | mc,c |
| VPB154-GW-082514-698-700 | WG | METHYL TERT-BUTYL ETHER | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | METHYLENE CHLORIDE | | 2.5 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | O-XYLENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | STYRENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | TETRACHLOROETHENE | 4.6 | 0.50 | UG/L | J | mc,s |
| VPB154-GW-082514-698-700 | WG | TOLUENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | TRANS-1,2-DICHLOROETHENE | | 0.50 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | TRANS-1,3-DICHLOROPROPENE | | 0.50 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-082514-698-700 | WG | TRICHLOROETHENE | 700 | 10 | UG/L | J | mc |
| VPB154-GW-082514-698-700 | WG | TRICHLOROFLUOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | VINYL CHLORIDE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-698-700 | WG | XYLENES, TOTAL | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,1,1-TRICHLOROETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,1,2,2-TETRACHLOROETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 10 | UG/L | UJ | mc,c |
| VPB154-GW-082514-718-720 | WG | 1,1,2-TRICHLOROETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,1-DICHLOROETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,1-DICHLOROETHENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2,4-TRICHLOROBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 15 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DIBROMOETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DICHLOROBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DICHLOROETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DICHLOROETHENE, TOTAL | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,2-DICHLOROPROPANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,3-DICHLOROBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 1,4-DICHLOROBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 2-BUTANONE | | 50 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 2-HEXANONE | | 50 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | 4-METHYL-2-PENTANONE | | 50 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-082514-718-720 | WG | ACETONE | | 50 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | BENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | BROMODICHLOROMETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | BROMOFORM | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | BROMOMETHANE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CARBON DISULFIDE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CARBON TETRACHLORIDE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CHLOROBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CHLOROETHANE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CHLOROFORM | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CHLOROMETHANE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CIS-1,2-DICHLOROETHENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CIS-1,3-DICHLOROPROPENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | CYCLOHEXANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | DIBROMOCHLOROMETHANE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | DICHLORODIFLUOROMETHANE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | ETHYLBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | ISOPROPYLBENZENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | M- AND P-XYLENE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | METHYL ACETATE | | 15 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | METHYL CYCLOHEXANE | | 10 | UG/L | UJ | mc,c |
| VPB154-GW-082514-718-720 | WG | METHYL TERT-BUTYL ETHER | | 10 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-082514-718-720 | WG | METHYLENE CHLORIDE | | 50 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | O-XYLENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | STYRENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TETRACHLOROETHENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TOLUENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TRANS-1,2-DICHLOROETHENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TRANS-1,3-DICHLOROPROPENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TRICHLOROETHENE | | 10 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | TRICHLOROFLUOROMETHANE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | VINYL CHLORIDE | | 20 | UG/L | UJ | mc |
| VPB154-GW-082514-718-720 | WG | XYLENES, TOTAL | | 30 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,1,1-TRICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,1,2,2-TETRACHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 1.0 | UG/L | UJ | mc,c |
| VPB154-GW-082514-738-740 | WG | 1,1,2-TRICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,1-DICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,1-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2,4-TRICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2-DIBROMOETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2-DICHLOROETHANE | | 1.0 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-082514-738-740 | WG | 1,2-DICHLOROETHENE, TOTAL | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,2-DICHLOROPROPANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,3-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 1,4-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 2-BUTANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 2-HEXANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | 4-METHYL-2-PENTANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | ACETONE | 15 | 5.0 | UG/L | J | mc,c |
| VPB154-GW-082514-738-740 | WG | BENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | BROMODICHLOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | BROMOFORM | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | BROMOMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CARBON DISULFIDE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CARBON TETRACHLORIDE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CHLOROBENZENE | 0.69 | 1.0 | UG/L | J | mc |
| VPB154-GW-082514-738-740 | WG | CHLOROETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CHLOROFORM | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CHLOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CIS-1,2-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CIS-1,3-DICHLOROPROPENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | CYCLOHEXANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | DIBROMOCHLOROMETHANE | | 1.0 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082514-738-740 | WG | DICHLORODIFLUOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | ETHYLBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | ISOPROPYLBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | M- AND P-XYLENE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | METHYL ACETATE | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | METHYL CYCLOHEXANE | | 1.0 | UG/L | UJ | mc,c |
| VPB154-GW-082514-738-740 | WG | METHYL TERT-BUTYL ETHER | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | METHYLENE CHLORIDE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | O-XYLENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | STYRENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | TETRACHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | TOLUENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | TRANS-1,2-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | TRANS-1,3-DICHLOROPROPENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | TRICHLOROETHENE | 1.4 | 1.0 | UG/L | J | mc |
| VPB154-GW-082514-738-740 | WG | TRICHLOROFUOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | VINYL CHLORIDE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082514-738-740 | WG | XYLENES, TOTAL | | 3.0 | UG/L | UJ | mc |
| VPB154-TRIP BLANK-082514 | WQ | 2-BUTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-082514 | WQ | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-082514 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | U | bt |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|------------|--------------------|-----|---------|
| WG148025-7 | 1,1-DICHLOROETHENE | 126 | 80-120% |
| | CARBON DISULFIDE | 129 | 80-120% |
| | ACETONE | 151 | 80-120% |
| | 2-BUTANONE | 125 | 80-120% |
| | 2-HEXANONE | 123 | 80-120% |

Associated samples: VPB154-GW-082214-663-665,VPB154-GW-082214-678-680,VPB154-GW-082514-698-700,VPB154-GW-082514-718-720,VPB154-GW-082514-738-740

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % D | Limits |
|------------|---------------------------------------|-----|--------|
| WG149034-4 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 39 | <20% |
| | METHYLCYCLOHEXANE | 29 | <20% |

Associated samples: VPB154-GW-082214-663-665,VPB154-GW-082214-678-680,VPB154-GW-082514-698-700,VPB154-GW-082514-718-720,VPB154-GW-082514-738-740

| | | | |
|------------|------------|-----|------|
| WG149439-4 | 2-BUTANONE | -21 | <20% |
| | 2-HEXANONE | -24 | <20% |

Associated samples: VPB154-TRIP BLANK-082514

Table A-3 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|--------------------------|------------------|--------|------|-------|--|
| VPB154-TRIP BLANK-082514 | CARBON DISULFIDE | 0.32 | 0.50 | UG/L | VPB154-GW-082214-663-665, VPB154-GW-082214-678-680, VPB154-GW-082514-698-700 |

Table A-4 - Surrogates

| Sample ID | Surrogate | % Recovery | Lower Limit | Upper Limit |
|--------------------------|-----------------------|------------|-------------|-------------|
| VPB154-GW-082514-698-700 | 1,2-DICHLOROETHANE-D4 | 122 | 70 | 120 |
| VPB154-GW-082214-663-665 | 1,2-DICHLOROETHANE-D4 | 121 | 70 | 120 |
| VPB154-GW-082514-718-720 | 1,2-DICHLOROETHANE-D4 | 122 | 70 | 120 |

Table A-5 - Matrix Spikes

| Sample ID | Compound | MS % Recovery | MSD % Recovery | Lower Limit | Upper Limit | RPD | RPD Limit |
|--------------------------|-------------|---------------|----------------|-------------|-------------|-----|-----------|
| VPB154-GW-082514-698-700 | CYCLOHEXANE | 148 | 147 | 71 | 133 | <1 | 30 |

| Sample ID | Compound | MS % Recovery | MSD % Recovery | Lower Limit | Upper Limit | RPD | RPD Limit |
|--------------------------|---------------------------------------|---------------|----------------|-------------|-------------|-----|-----------|
| VPB154-GW-082514-698-700 | O-XYLENE | 120 | 122 | 80 | 120 | 1 | 30 |
| VPB154-GW-082514-698-700 | XYLEMES, TOTAL | 121 | 122 | 89 | 116 | 0 | 30 |
| VPB154-GW-082514-698-700 | CIS-1,2-DICHLOROETHENE | 125 | 129 | 70 | 125 | 3 | 30 |
| VPB154-GW-082514-698-700 | METHYL TERT-BUTYL ETHER | 134 | 133 | 65 | 125 | 1 | 30 |
| VPB154-GW-082514-698-700 | 1,2-DICHLOROETHENE, TOTAL | 126 | 130 | 84 | 121 | 3 | 30 |
| VPB154-GW-082514-698-700 | 2-HEXANONE | 150 | 162 | 55 | 130 | 8 | 30 |
| VPB154-GW-082514-698-700 | BENZENE | 119 | 121 | 80 | 120 | 2 | 30 |
| VPB154-GW-082514-698-700 | 1,1,1-TRICHLOROETHANE | 135 | 131 | 65 | 130 | 3 | 30 |
| VPB154-GW-082514-698-700 | 4-METHYL-2-PENTANONE | 139 | 132 | 60 | 135 | 5 | 30 |
| VPB154-GW-082514-698-700 | METHYL CYCLOHEXANE | 174 | 173 | 73 | 125 | 1 | 30 |
| VPB154-GW-082514-698-700 | BROMODICHLOROMETHANE | 127 | 123 | 75 | 120 | 3 | 30 |
| VPB154-GW-082514-698-700 | 1,1-DICHLOROETHANE | 138 | 135 | 70 | 135 | 2 | 30 |
| VPB154-GW-082514-698-700 | 1,1-DICHLOROETHENE | 141 | 146 | 70 | 130 | 3 | 30 |
| VPB154-GW-082514-698-700 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 153 | 165 | 73 | 126 | 6 | 30 |
| VPB154-GW-082514-698-700 | 1,2-DICLOROPROPANE | 126 | 127 | 75 | 125 | 0 | 30 |
| VPB154-GW-082514-698-700 | TRICHLOROETHENE | 0 | 0 | 70 | 125 | 5 | 30 |
| VPB154-GW-082514-698-700 | METHYL ACETATE | 140 | 132 | 70 | 132 | 6 | 30 |

Table A-6 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|---------------------------------------|----------------|-------------|-------------|--|
| WG149034-1 | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 139 | 73 | 126 | VPB154-GW-082214-663-665 VPB154-GW-082214-678-680 VPB154-GW-082514-698-700 |

Attachment B
Qualifier Codes and Explanations

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|--------------------|--|
| be | Equipment blank contamination |
| bt | Trip blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |
| mc | Method compliance nonconformance |



600 Technology Way
Scarborough, ME 04074
Tel: (207) 874-2400
Fax: (207) 775-4029

CHAIN of CUSTODY

**PLEASE BEAR DOWN AND
PRINT LEGIBLY IN PEN**

Page _____ of _____

| | | | | | | | | | | | | |
|---|-----------------------------------|--|---------------------------------|--|--------------------------|------------------------------|----------------|--------------------------|----------------|----------------|----------------|----------------|
| Client Resolution Consultants | | Contact T. Viverda | Phone # (315)475-1980 | Fax # () | | | | | | | | |
| Address 100 Red Schoolhouse Rd. | | City Chestertown | State NY | Zip Code 10977 | | | | | | | | |
| Purchase Order # | | Proj. Name / No. MWAP Deltapage/60066526 | Katahdin Quote # | | | | | | | | | |
| Bill (if different than above) | | Address | | | | | | | | | | |
| Sampler (Print / Sign) Vincent Varrichio N/A | | | | | Copies To: | | | | | | | |
| LAB USE ONLY | | WORK ORDER #: SH 6880 | | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | |
| | | KATAHDIN PROJECT NUMBER | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON |
| REMARKS: | | | | | | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | | | |
| TEMP°C _____ | | <input type="checkbox"/> TEMP BLANK | <input type="checkbox"/> INTACT | <input type="checkbox"/> NOT INTACT | | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | | | | | | | | |
| | UPD154-GW-082514-663665 | 8/25/14 / 1120 | GW | 3 | 3 | | | | | | | |
| | UPD154-GW-082514-C781610 | 8/25/14 / 1325 | GW | 3 | 3 | | | | | | | |
| | UPD154-GW-082514-698-70 | 8/25/14 / 1010 | GW | 3 | 3 | | | | | | | |
| | UPD154-GW-MS MSD - 082514-698-700 | 8/25/14 / 1010 | GW | 6 | 6 | | | | | | | |
| | UPD154-GW-082514-718-720 | 8/25/14 / 1340 | GW | 3 | 3 | | | | | | | |
| | UPD154-Trip Blank-082514 | 8/25/14 / 1430 | W | 3 | 3 | | | | | | | |
| | UPD154-GW-082514-738-740 | 8/25/14 / 1515 | GW | 3 | 3 | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| | | / | | | | | | | | | | |
| COMMENTS UPD154-GW-082514-718-720 - Very Silty also UPD154-GW-082514-738-740 | | | | | | | | | | | | |
| Relinquished By: (Signature) | Date / Time 8/25/14 1610 | Received By: (Signature) FedEx | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | | | | |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | | | | |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

KAS-COC1

Dengelo

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6880-1
 Client ID: 154-082214-663-665
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6880
 Lab File ID: C8765.D

Sample Date: 22-AUG-14
 Received Date: 26-AUG-14
 Extract Date: 27-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149034

Analysis Date: 27-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | + T | 0.26 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | T | 16 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + U T | 4.29 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | + T | 84 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | T | 10 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | T | 2.6 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | T | 3.9 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | T | 1.6 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | T | 2.0 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | J T | 2.4 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | T | 2.4 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | X | 620 600 | ug/L | X 10 | 1 | 1.0 | 0.28 2.8 | 0.56 5.0 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | T | 1.2 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Riz 12/29/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-1
Client ID: 154-082214-663-665
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8765.D

Sample Date: 22-AUG-14 **Analysis Date:** 27-AUG-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 27-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149034 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | T | 3.9 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 86.8 | % | | | | | |
| Toluene-d8 | | 91.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 121. | % | | | | | |
| Dibromofluoromethane | | 107. | % | | | | | |

12/29/14

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6880-2
 Client ID: 154-082214-678-680
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6880
 Lab File ID: C8766.D

Sample Date: 22-AUG-14 Analysis Date: 27-AUG-14
 Received Date: 26-AUG-14 Analyst: REC
 Extract Date: 27-AUG-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG149034 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|----------|----------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | J | 3.2 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | + U J | 0.28 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | + J | 21 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | J | 8.7 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | J | 0.97 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | | 2.1 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | J | 0.78 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | J | 0.37 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | P | 580 SSD | ug/L | X(O) | 1 | 1010 | 0.28 2.8 | 0.50 S-O |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | J | 0.53 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 2.8 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | J | 0.39 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2



Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-2
Client ID: 154-082214-678-680
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8766.D

Sample Date: 22-AUG-14 **Analysis Date:** 27-AUG-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 27-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149034 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|-------------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | | 2.1 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 84.0 | % | | | | | |
| Toluene-d8 | | 88.4 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 120. | % | | | | | |
| Dibromofluoromethane | | 106. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6880-3
 Client ID: 154-082514-698-700
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6880
 Lab File ID: C8767.D

Sample Date: 25-AUG-14
 Received Date: 26-AUG-14
 Extract Date: 27-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149034

Analysis Date: 27-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 0.5 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | MM | 0.5 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.33 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | LMM | 0.30 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 0.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | | 26 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | UMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | MM | 1.5 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | MM | 2.9 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | | 1.2 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | JMM | 0.60 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | | 5.3 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | UMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | UM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | EMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,2-Dichloropropane | UMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | UMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | UM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | UM | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | | 0.68 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | | 4.6 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | UMM | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-3
Client ID: 154-082514-698-700
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8767.D

Sample Date: 25-AUG-14 **Analysis Date:** 27-AUG-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 27-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149034 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | UMM | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | + U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | UMM | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | ULMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | UMM | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | MM | 2.9 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 87.9 | % | | | | | |
| Toluene-d8 | | 89.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 122. | % | | | | | |
| Dibromofluoromethane | | 106. | % | | | | | |

12/29/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-4DL
Client ID: 154-082514-718-720
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8763.D

Sample Date: 25-AUG-14 **Analysis Date:** 27-AUG-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 27-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149034 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 20 | ug/L | 20 | 2 | 40. | 4.8 | 20. |
| Chloromethane | U | 20 | ug/L | 20 | 2 | 40. | 7.2 | 20. |
| Vinyl Chloride | U | 20 | ug/L | 20 | 2 | 40. | 5.0 | 20. |
| Bromomethane | U | 20 | ug/L | 20 | 2 | 40. | 9.8 | 20. |
| Chloroethane | U | 20 | ug/L | 20 | 2 | 40. | 11. | 20. |
| Trichlorodifluoromethane | U | 20 | ug/L | 20 | 2 | 40. | 4.8 | 20. |
| 1,1-Dichloroethene | U | 10 | ug/L | 20 | 1 | 20. | 7.0 | 10. |
| Carbon Disulfide | U | 10 | ug/L | 20 | 1 | 20. | 5.0 | 10. |
| Freon-113 | UL | 10 | ug/L | 20 | 1 | 20. | 6.2 | 10. |
| Methylene Chloride | U | 50 | ug/L | 20 | 5 | 100 | 23. | 50. |
| Acetone | U | 50 | ug/L | 20 | 5 | 100 | 44. | 50. |
| trans-1,2-Dichloroethene | U | 10 | ug/L | 20 | 1 | 20. | 5.0 | 10. |
| Methyl tert-butyl Ether | U | 10 | ug/L | 20 | 1 | 20. | 7.2 | 10. |
| 1,1-Dichloroethane | U | 10 | ug/L | 20 | 1 | 20. | 4.2 | 10. |
| cis-1,2-Dichloroethene | U | 10 | ug/L | 20 | 1 | 20. | 4.2 | 10. |
| Chloroform | U | 10 | ug/L | 20 | 1 | 20. | 6.4 | 10. |
| 1,1,1-Trichloroethane | U | 10 | ug/L | 20 | 1 | 20. | 4.0 | 10. |
| 2-Butanone | U | 50 | ug/L | 20 | 5 | 100 | 26. | 50. |
| Cyclohexane | U | 10 | ug/L | 20 | 1 | 20. | 6.2 | 10. |
| Carbon Tetrachloride | U | 10 | ug/L | 20 | 1 | 20. | 4.4 | 10. |
| Benzene | U | 10 | ug/L | 20 | 1 | 20. | 5.2 | 10. |
| 1,2-Dichloroethane | U | 10 | ug/L | 20 | 1 | 20. | 4.0 | 10. |
| Trichloroethene | U | 10 | ug/L | 20 | 1 | 20. | 5.6 | 10. |
| 1,2-Dichloropropane | U | 10 | ug/L | 20 | 1 | 20. | 5.0 | 10. |
| Bromodichloromethane | U | 10 | ug/L | 20 | 1 | 20. | 6.6 | 10. |
| cis-1,3-Dichloropropene | U | 10 | ug/L | 20 | 1 | 20. | 3.8 | 10. |
| Toluene | U | 10 | ug/L | 20 | 1 | 20. | 5.4 | 10. |
| 4-Methyl-2-Pentanone | U | 50 | ug/L | 20 | 5 | 100 | 26. | 50. |
| trans-1,3-Dichloropropene | U | 10 | ug/L | 20 | 1 | 20. | 4.0 | 10. |
| 1,1,2-Trichloroethane | U | 10 | ug/L | 20 | 1 | 20. | 6.6 | 10. |
| Tetrachloroethene | U | 10 | ug/L | 20 | 1 | 20. | 8.0 | 10. |
| Dibromochloromethane | U | 10 | ug/L | 20 | 1 | 20. | 6.0 | 10. |
| 2-Hexanone | U | 50 | ug/L | 20 | 5 | 100 | 34. | 50. |
| Chlorobenzene | U | 10 | ug/L | 20 | 1 | 20. | 4.4 | 10. |
| Ethylbenzene | U | 10 | ug/L | 20 | 1 | 20. | 4.2 | 10. |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH6880-4DL
 Client ID: 154-082514-718-720
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6880
 Lab File ID: C8763.D

Sample Date: 25-AUG-14 Analysis Date: 27-AUG-14
 Received Date: 26-AUG-14 Analyst: REC
 Extract Date: 27-AUG-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG149034 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 30 | ug/L | 20 | 3 | 60. | 5.0 | 30. |
| Styrene | U | 10 | ug/L | 20 | 1 | 20. | 4.6 | 10. |
| Bromoform | U | 10 | ug/L | 20 | 1 | 20. | 4.6 | 10. |
| Isopropylbenzene | U | 10 | ug/L | 20 | 1 | 20. | 4.6 | 10. |
| 1,1,2,2-Tetrachloroethane | U | 10 | ug/L | 20 | 1 | 20. | 7.6 | 10. |
| 1,3-Dichlorobenzene | U | 10 | ug/L | 20 | 1 | 20. | 5.2 | 10. |
| 1,4-Dichlorobenzene | U | 10 | ug/L | 20 | 1 | 20. | 4.8 | 10. |
| 1,2-Dichlorobenzene | U | 10 | ug/L | 20 | 1 | 20. | 3.0 | 10. |
| 1,2,4-Trichlorobenzene | U | 10 | ug/L | 20 | 1 | 20. | 7.4 | 10. |
| Methyl Acetate | U | 15 | ug/L | 20 | 1 | 20. | 11. | 15. |
| Methylcyclohexane | UL | 10 | ug/L | 20 | 1 | 20. | 6.0 | 10. |
| o-Xylene | U | 10 | ug/L | 20 | 1 | 20. | 5.0 | 10. |
| M+P-Xylenes | U | 20 | ug/L | 20 | 2 | 40. | 12. | 20. |
| 1,2-Dichloroethylene (Total) | U | 20 | ug/L | 20 | 2 | 40. | 4.2 | 20. |
| 1,2-Dibromoethane | U | 10 | ug/L | 20 | 1 | 20. | 4.4 | 10. |
| 1,2-Dibromo-3-Chloropropane | U | 15 | ug/L | 20 | 1 | 20. | 10. | 15. |
| P-Bromofluorobenzene | | 89.2 | % | | | | | |
| Toluene-d8 | | 94.6 | % | | | | | |
| 1,2-Dichloroethane-d4 | * | 122. | % | | | | | |
| Dibromofluoromethane | | 107. | % | | | | | |



Report of Analytical Results

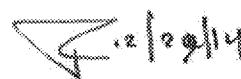
Client: ENSAFE
 Lab ID: SH6880-6DL
 Client ID: 154-082514-738-740
 Project: Navy Clean WE15-03-06 NW
 SDG: SH6880
 Lab File ID: C8764.D

Sample Date: 25-AUG-14
 Received Date: 26-AUG-14
 Extract Date: 27-AUG-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149034

Analysis Date: 27-AUG-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD | |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|-----|
| Dichlorodifluoromethane | U | US | 2.0 | ug/L | 2 | 2 | 4.0 | 0.48 | 2.0 |
| Chloromethane | U | | 2.0 | ug/L | 2 | 2 | 4.0 | 0.72 | 2.0 |
| Vinyl Chloride | U | | 2.0 | ug/L | 2 | 2 | 4.0 | 0.50 | 2.0 |
| Bromomethane | U | | 2.0 | ug/L | 2 | 2 | 4.0 | 0.98 | 2.0 |
| Chloroethane | U | | 2.0 | ug/L | 2 | 2 | 4.0 | 1.1 | 2.0 |
| Trichlorofluoromethane | U | | 2.0 | ug/L | 2 | 2 | 4.0 | 0.48 | 2.0 |
| 1,1-Dichloroethene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.70 | 1.0 |
| Carbon Disulfide | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Freon-113 | UL | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.62 | 1.0 |
| Methylene Chloride | U | | 5.0 | ug/L | 2 | 5 | 10. | 2.3 | 5.0 |
| Acetone | J | 15 | ug/L | 2 | 5 | 10. | 4.4 | 5.0 | |
| trans-1,2-Dichloroethene | U | US | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Methyl tert-butyl Ether | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.72 | 1.0 |
| 1,1-Dichloroethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |
| cis-1,2-Dichloroethene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |
| Chloroform | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.64 | 1.0 |
| 1,1,1-Trichloroethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| 2-Butanone | U | | 5.0 | ug/L | 2 | 5 | 10. | 2.6 | 5.0 |
| Cyclohexane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.62 | 1.0 |
| Carbon Tetrachloride | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 |
| Benzene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.52 | 1.0 |
| 1,2-Dichloroethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| Trichloroethene | J | 1.4 | ug/L | 2 | 1 | 2.0 | 0.56 | 1.0 | |
| 1,2-Dichloropropane | U | US | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Bromodichloromethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.66 | 1.0 |
| cis-1,3-Dichloropropene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.38 | 1.0 |
| Toluene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.54 | 1.0 |
| 4-Methyl-2-Pentanone | U | | 5.0 | ug/L | 2 | 5 | 10. | 2.6 | 5.0 |
| trans-1,3-Dichloropropene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| 1,1,2-Trichloroethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.66 | 1.0 |
| Tetrachloroethene | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.80 | 1.0 |
| Dibromochloromethane | U | | 1.0 | ug/L | 2 | 1 | 2.0 | 0.60 | 1.0 |
| 2-Hexanone | U | | 5.0 | ug/L | 2 | 5 | 10. | 3.4 | 5.0 |
| Chlorobenzene | J | 0.69 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 | |
| Ethylbenzene | U | US | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-6DL
Client ID: 154-082514-738-740
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8764.D

Sample Date: 25-AUG-14 **Analysis Date:** 27-AUG-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 27-AUG-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149034 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 3.0 | ug/L | 2 | 3 | 6.0 | 0.50 | 3.0 |
| Styrene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| Bromoform | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| Isopropylbenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| 1,1,2,2-Tetrachloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.76 | 1.0 |
| 1,3-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.52 | 1.0 |
| 1,4-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.48 | 1.0 |
| 1,2-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.30 | 1.0 |
| 1,2,4-Trichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.74 | 1.0 |
| Methyl Acetate | U | 1.5 | ug/L | 2 | 1 | 2.0 | 1.1 | 1.5 |
| Methylcyclohexane | UL | 1.0 | ug/L | 2 | 1 | 2.0 | 0.60 | 1.0 |
| o-Xylene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| M+P-Xylenes | U | 2.0 | ug/L | 2 | 2 | 4.0 | 1.2 | 2.0 |
| 1,2-Dichloroethylene (Total) | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.42 | 2.0 |
| 1,2-Dibromoethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | U | 1.5 | ug/L | 2 | 1 | 2.0 | 1.0 | 1.5 |
| P-Bromofluorobenzene | | 90.4 | % | | | | | |
| Toluene-d8 | | 94.6 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 119. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |



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Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-5RA
Client ID: VPB154-TB-082514
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8841.D

Sample Date: 25-AUG-14
Received Date: 26-AUG-14
Extract Date: 03-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | J | 0.53 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | -U- U | 0.32 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | UL | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | -U- UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | -U- UJ | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0* | 0.21 | 0.50 |

Page 1 of 2



Report of Analytical Results

Client: ENSAFE
Lab ID: SH6880-5RA
Client ID: VPB154-TB-082514
Project: Navy Clean WE15-03-06 NW
SDG: SH6880
Lab File ID: C8841.D

Sample Date: 25-AUG-14 **Analysis Date:** 03-SEP-14
Received Date: 26-AUG-14 **Analyst:** REC
Extract Date: 03-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149439 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.9 | % | | | | | |
| Toluene-d8 | | 94.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 104. | % | | | | | |
| Dibromofluoromethane | | 93.0 | % | | | | | |



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Data Validation Report

| | | |
|-----------------------|---|--------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH7085 | |
| Analyses/Method: | EPA SW-846 Method 8260B for VOCs (GC/MS) | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/05/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH7085_8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 27 and 28, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-GW-082714-763-765 | Groundwater |
| VPB154-GW-082714-778-780 | Groundwater |
| VPB154-GW-082814-798-800 | Groundwater |
| VPB154-GW-082814-818-820 | Groundwater |
| VPB154-TRIP BLANK-082814 | Trip Blank |

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- Data completeness (chain-of-custody [COC])/sample integrity
- Holding times and sample preservation
- GC/MS performance checks
- Initial calibration/continuing calibration verification
- Laboratory blanks/equipment blanks/trip blanks
- Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- Laboratory control sample (LCS) results

- NA Field duplicate results
✓ Internal standard results
✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

The vials of sample VPB154-GW-082714-778-780 were mostly soil and had very little standing water. Therefore, each vial was decanted and composited into one vial and analyzed. Due to limited sample volume, the sample was analyzed at a dilution of 1:2. Positive and nondetect results for this sample were qualified as estimated (J and UJ, respectively), due to possible loss of sample integrity during the decanting procedure. Qualified sample results are shown in Table 1.

Holding Times and Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Tables A-1 and A-2.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

ICV Recovery Nonconformances:

| Nonconformance | Actions | |
|---------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R > 120% | J | No qualification |
| 20% < %R < 80% | J | UJ |
| %R < 20% (see note) | J | R* |

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject (R) sample results previously negated (U) on the basis of blank contamination.

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift > 20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Tables A-3 and A-4.

Sample results were qualified as follows:

| Blank type | Blank result | Sample result | Action for samples |
|--|---------------------------|---|---|
| Method, Storage, Field, Trip, or Instrument* | Detects \leq LOQ | Not detected | No qualification |
| | | < LOQ | Report sample LOQ value with a U |
| | | \geq LOQ and \leq 2x LOQ | Report the sample result with a U** |
| | > LOQ | \geq 2x the LOQ | No qualifications |
| | | < LOQ | Report sample LOQ value with a U |
| | | \geq LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R |
| | | \geq LOQ and \geq blank contamination | If the result is \leq 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.** |

* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.

**Based on professional judgment.

LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria.

Nonconformances are summarized in Attachment A in Table A-5.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|--------------------------------------|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) | J | R |
| (LL = lower limit, UL = upper limit) | | |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082714-763-765 | WG | ACETONE | 13 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082714-763-765 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |
| VPB154-GW-082714-778-780 | WG | 1,1,1-TRICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,1,2,2-TETRACHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,1,2-TRICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,1-DICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,1-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2,4-TRICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DIBROMOETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DICHLOROETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DICHLOROETHENE, TOTAL | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,2-DICHLOROPROPANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,3-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 1,4-DICHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 2-BUTANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 2-HEXANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | 4-METHYL-2-PENTANONE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | ACETONE | 38 | 5.0 | UG/L | J | mc,c,l |
| VPB154-GW-082714-778-780 | WG | BENZENE | | 1.0 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082714-778-780 | WG | BROMODICHLOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | BROMOFORM | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | BROMOMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CARBON DISULFIDE | | 2.0* | UG/L | UJ | mc,bl |
| VPB154-GW-082714-778-780 | WG | CARBON TETRACHLORIDE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CHLOROBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CHLOROETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CHLOROFORM | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CHLOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CIS-1,2-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CIS-1,3-DICHLOROPROPENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | CYCLOHEXANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | DIBROMOCHLOROMETHANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | DICHLORODIFLUOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | ETHYLBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | ISOPROPYLBENZENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | M- AND P-XYLENE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | METHYL ACETATE | | 1.5 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | METHYL CYCLOHEXANE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | METHYL TERT-BUTYL ETHER | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | METHYLENE CHLORIDE | | 5.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | O-XYLENE | | 1.0 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-082714-778-780 | WG | STYRENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TETRACHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TOLUENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TRANS-1,2-DICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TRANS-1,3-DICHLOROPROPENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TRICHLOROETHENE | | 1.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | TRICHLOROFLUOROMETHANE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | VINYL CHLORIDE | | 2.0 | UG/L | UJ | mc |
| VPB154-GW-082714-778-780 | WG | XYLEMES, TOTAL | | 3.0 | UG/L | UJ | mc |
| VPB154-GW-082814-798-800 | WG | 2-BUTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082814-798-800 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082814-798-800 | WG | ACETONE | 13 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082814-798-800 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bt |
| VPB154-GW-082814-818-820 | WG | 2-BUTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082814-818-820 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082814-818-820 | WG | ACETONE | 5.9 | 2.5 | UG/L | J | c,l |
| VPB154-GW-082814-818-820 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | bt |
| VPB154-TRIP BLANK-082814 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 - Initial Calibration Verification Standard**

| ICV ID | Compound | % R | Limits |
|---|----------|-----|---------|
| WG149371-7 | ACETONE | 173 | 80-120% |
| Associated samples: all samples in SDG SH7085 | | | |

Table A-2 -Continuing Calibration Verification Standard

| CCV ID | Compound | % D | Limits |
|--|------------|-----|--------|
| WG149439-4 | 2-BUTANONE | -21 | <20% |
| | 2-HEXANONE | -24 | <20% |
| Associated samples: VPB154-GW-082814-798-800, VPB154-GW-082814-818-820 | | | |

Table A-3 - Lab Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------|------------------|--------|------|-------|--|
| WG149371-9 | CARBON DISULFIDE | 0.40 | 0.50 | UG/L | VPB154-GW-082714-763-765 VPB154-GW-082714-778-780 VPB154-TRIP BLANK-082814 |

Table A-4 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|--------------------------|------------------|--------|------|-------|---|
| VPB154-TRIP BLANK-082814 | CARBON DISULFIDE | 0.35 | 0.50 | UG/L | VPB154-GW-082814-798-800, VPB154-GW-082814-818-820 |

Table A-5 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|----------|----------------|-------------|-------------|--|
| WG149439-1 | ACETONE | 146 | 40 | 140 | VPB154-GW-082814-798-800 VPB154-GW-082814-818-820 |
| WG149371-8 | ACETONE | 173 | 40 | 140 | VPB154-GW-082714-763-765 VPB154-GW-082714-778-780 |

Attachment B
Qualifier Codes and Explanations

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bt | Trip blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |
| mc | Method compliance nonconformance |



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CHAIN of CUSTODY

PLEASE BEAR DOWN AND
PRINT LEGIBLY IN PEN

Page 1 of 1

| | | | | | | | | | | | |
|--|---|---|---|------------------------------------|--|---------------|---------------|---------------|---------------|---------------|---------------|
| Client <i>Resolution (Env) Inc.</i> | Contact <i>E. Virelizier</i> | Phone # <i>(415) 405-4110</i> | Fax # <i>()</i> | | | | | | | | |
| Address <i>160 Rte Schoolhouse Rd.</i> | City <i>Chestnut Ridge</i> | State <i>NY</i> | Zip Code <i>10977</i> | | | | | | | | |
| Purchase Order # | Proj. Name / No. <i>WV1010 Deltabarge / 6626CJ26</i> | Katahdin Quote # | | | | | | | | | |
| Bill (if different than above) | Address | | | | | | | | | | |
| Sampler (Print / Sign) <i>Vincent Virelizier / V.V.</i> | | | Copies To: | | | | | | | | |
| LAB USE ONLY | WORK ORDER #: <i>SH 7085</i> | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | | |
| KATAHDIN PROJECT NUMBER _____ | | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON | Fill OY ON |
| REMARKS: | | | | | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | | |
| TEMP°C _____ <input type="checkbox"/> TEMP BLANK <input type="checkbox"/> INTACT <input type="checkbox"/> NOT INTACT | | | | | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | 1 | 2 | 3 | 4 | 5 | 6 | 7 |
| | <i>VPD154-6L-082814-763-705</i> | <i>8/28/14/1210</i> | <i>6W</i> | <i>3</i> | <i>1</i> | | | | | | |
| | <i>VPD154-6L-082814-778-780</i> | <i>8/28/14/1415</i> | <i>6W</i> | <i>3</i> | <i>1</i> | | | | | | |
| | <i>VPD154-6L-082814-798-802</i> | <i>8/28/14/1045</i> | <i>6W</i> | <i>3</i> | <i>1</i> | | | | | | |
| | <i>VPD154-6L-082814-818-812</i> | <i>8/28/14/1140</i> | <i>6W</i> | <i>3</i> | <i>1</i> | | | | | | |
| | <i>VPD154-Trip1BLK-082814</i> | <i>8/28/14/1450</i> | <i>W</i> | <i>3</i> | <i>1</i> | | | | | | |
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| COMMENTS | | | | | | | | | | | |
| Relinquished By: (Signature) <i>V.V.</i> | Date / Time <i>8/28/14/1600</i> | Received By: (Signature) <i>Mary Zabel</i> | Relinquished By: (Signature) <i>Mary Zabel</i> | Date / Time <i>9/29/14/1600</i> | Received By: (Signature) <i>Z. 8-29-14 0930</i> | | | | | | |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | | | | | | |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN
SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7085-1
 Client ID: 154-082714-763-765
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7085
 Lab File ID: C8824.D

Sample Date: 27-AUG-14 Analysis Date: 02-SEP-14
 Received Date: 29-AUG-14 Analyst: REC
 Extract Date: 02-SEP-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG149371 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------------------|----------------------------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U U | 0.29 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U T | 13 13 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-1
Client ID: 154-082714-763-765
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8824.D

Sample Date: 27-AUG-14
Received Date: 29-AUG-14
Extract Date: 02-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149371

Analysis Date: 02-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 90.8 | % | | | | | |
| Toluene-d8 | | 91.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 119. | % | | | | | |
| Dibromofluoromethane | | 104. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7085-2DL
 Client ID: 154-082714-778-780
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7085
 Lab File ID: C8823.D

Sample Date: 27-AUG-14
 Received Date: 29-AUG-14
 Extract Date: 02-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149371

Analysis Date: 02-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|-------------|-------------|----------|----------|------------|-------------|------------|
| Dichlorodifluoromethane | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.48 | 2.0 |
| Chloromethane | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.72 | 2.0 |
| Vinyl Chloride | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.50 | 2.0 |
| Bromomethane | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.98 | 2.0 |
| Chloroethane | U | 2.0 | ug/L | 2 | 2 | 4.0 | 1.1 | 2.0 |
| Trichlorofluoromethane | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.48 | 2.0 |
| 1,1-Dichloroethene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.70 | 1.0 |
| Carbon Disulfide | | 0.57 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Freon-113 | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.62 | 1.0 |
| Methylene Chloride | U | 5.0 | ug/L | 2 | 5 | 10. | 2.3 | 5.0 |
| Acetone | U | 38 | ug/L | 2 | 5 | 10. | 4.4 | 5.0 |
| trans-1,2-Dichloroethene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Methyl tert-butyl Ether | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.72 | 1.0 |
| 1,1-Dichloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |
| cis-1,2-Dichloroethene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |
| Chloroform | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.64 | 1.0 |
| 1,1,1-Trichloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| 2-Butanone | U | 5.0 | ug/L | 2 | 5 | 10. | 2.6 | 5.0 |
| Cyclohexane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.62 | 1.0 |
| Carbon Tetrachloride | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 |
| Benzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.52 | 1.0 |
| 1,2-Dichloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| Trichloroethene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.56 | 1.0 |
| 1,2-Dichloropropane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| Bromodichloromethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.66 | 1.0 |
| cis-1,3-Dichloropropene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.38 | 1.0 |
| Toluene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.54 | 1.0 |
| 4-Methyl-2-Pentanone | U | 5.0 | ug/L | 2 | 5 | 10. | 2.6 | 5.0 |
| trans-1,3-Dichloropropene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.40 | 1.0 |
| 1,1,2-Trichloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.66 | 1.0 |
| Tetrachloroethene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.80 | 1.0 |
| Dibromochloromethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.60 | 1.0 |
| 2-Hexanone | U | 5.0 | ug/L | 2 | 5 | 10. | 3.4 | 5.0 |
| Chlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 |
| Ethylbenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.42 | 1.0 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-2DL
Client ID: 154-082714-778-780
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8823.D

Sample Date: 27-AUG-14 **Analysis Date:** 02-SEP-14
Received Date: 29-AUG-14 **Analyst:** REC
Extract Date: 02-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149371 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | VS | ug/L | 2 | 3 | 6.0 | 0.50 | 3.0 |
| Styrene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| Bromoform | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| Isopropylbenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.46 | 1.0 |
| 1,1,2,2-Tetrachloroethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.76 | 1.0 |
| 1,3-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.52 | 1.0 |
| 1,4-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.48 | 1.0 |
| 1,2-Dichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.30 | 1.0 |
| 1,2,4-Trichlorobenzene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.74 | 1.0 |
| Methyl Acetate | U | 1.5 | ug/L | 2 | 1 | 2.0 | 1.1 | 1.5 |
| Methylcyclohexane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.60 | 1.0 |
| o-Xylene | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.50 | 1.0 |
| M+P-Xylenes | U | 2.0 | ug/L | 2 | 2 | 4.0 | 1.2 | 2.0 |
| 1,2-Dichloroethylene (Total) | U | 2.0 | ug/L | 2 | 2 | 4.0 | 0.42 | 2.0 |
| 1,2-Dibromoethane | U | 1.0 | ug/L | 2 | 1 | 2.0 | 0.44 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | U | 1.5 | ug/L | 2 | 1 | 2.0 | 1.0 | 1.5 |
| P-Bromofluorobenzene | | 90.6 | % | | | | | |
| Toluene-d8 | | 91.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 118. | % | | | | | |
| Dibromofluoromethane | | 102. | % | | | | | |



Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-3RA
Client ID: 154-082814-798-800
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8845.D

Sample Date: 28-AUG-14
Received Date: 29-AUG-14
Extract Date: 03-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|--------------|-----------------|-----------------|--------------|--------------|----------------|-----------------|-----------------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.26 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 13 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-3RA
Client ID: 154-082814-798-800
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8845.D

Sample Date: 28-AUG-14 **Analysis Date:** 03-SEP-14
Received Date: 29-AUG-14 **Analyst:** REC
Extract Date: 03-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149439 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.9 | % | | | | | |
| Toluene-d8 | | 93.9 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 107. | % | | | | | |
| Dibromofluoromethane | | 95.0 | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7085-4RA
 Client ID: 154-082814-818-820
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7085
 Lab File ID: C8846.D

Sample Date: 28-AUG-14
 Received Date: 29-AUG-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|--|--------------------------------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U S | 0.29 1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U T | 5.9 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U U T | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U U T | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-4RA
Client ID: 154-082814-818-820
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8846.D

Sample Date: 28-AUG-14
Received Date: 29-AUG-14
Extract Date: 03-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 94.1 | % | | | | | |
| Toluene-d8 | | 94.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 110. | % | | | | | |
| Dibromofluoromethane | | 94.8 | % | | | | | |

Report of Analytical Results

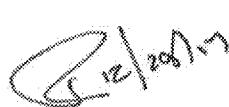
Client: ENSAFE
Lab ID: SH7085-5
Client ID: VPB154-TB-082814
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8818.D

Sample Date: 28-AUG-14
Received Date: 29-AUG-14
Extract Date: 02-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149371

Analysis Date: 02-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | J | 0.58 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- | 0.35 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | J | 1.4 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | UL | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7085-5
Client ID: VPB154-TB-082814
Project: Navy Clean WE15-03-06 NW
SDG: SH7085
Lab File ID: C8818.D

Sample Date: 28-AUG-14
Received Date: 29-AUG-14
Extract Date: 02-SEP-14
Extracted By: REC
Extraction Method: SW846 5030
Lab Prep Batch: WG149371

Analysis Date: 02-SEP-14
Analyst: REC
Analysis Method: SW846 8260C
Matrix: AQ
% Solids: NA
Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 87.8 | % | | | | | |
| Toluene-d8 | | 90.0 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 100. | % | | | | | |
| Dibromofluoromethane | | 98.2 | % | | | | | |



Resolution Consultants
250 Apollo Drive
Chelmsford, MA 01824

978.905.2100 tel
978.905.2101 fax

Data Validation Report

| | | |
|--|---|-----------------------------------|
| Project: | Regional Groundwater Investigation - NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Service Request: | SH7193 | |
| Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS) and Standard Method 5310 for Total Organic Carbon by High-Temperature Combustion | | |
| Validation Level: | 3 | |
| AECOM Project Number: | 60266526.SA.DV | |
| Prepared by: | Dawn Brule/RESCON | Completed on: 12/19/2014 |
| Reviewed by: | Lori Herberich/RESCON | File Name: SH7193_5310B and 8260B |

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 29, 2014 and September 2, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-EB-090214 | Equipment blank |
| VPB154-GW-082914-838-840 | Groundwater |
| VPB154-GW-082914-858-860 | Groundwater |
| VPB154-GW-090214-908-910 | Groundwater |
| VPB154-TRIP BLANK-090214 | Trip Blank |

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Standard Methods for the Examination of Water and Wastewater, Method SM5310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and *Quality Systems Manual (QSM)* for Environmental Laboratories, Version 4.2 (DoD, October 2010) where applicable. In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- Data completeness (chain-of-custody [COC])/sample integrity

- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✗ Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✗ Laboratory control sample (LCS)/laboratory control sample duplicate (LCSD) results
- NA Field duplicate results
- ✓ Internal standard results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

The vials of sample VPB154-GW-090214-908-910 were mostly soil and had very little standing water. Therefore, the laboratory decanted the water from the individual vials into one vial as a composite. Due to limited sample volume, the sample was analyzed at a 1:50 dilution. Positive and nondetect results for this sample were qualified as estimated (J and UJ, respectively), due to possible loss of sample integrity during the decanting procedure.

Qualified sample results are shown in Table 1.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Table A-1.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

CCV Linearity Nonconformances:

| Nonconformance | Actions | |
|----------------|------------------|---------------------|
| | Detected Results | Nondetected Results |
| %D > 20% | J | UJ |
| %Drift >20% | J* | UJ* |

* No guidance in NFG, thus professional judgment was used

Qualified sample results are shown in Table 1.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL).

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method, equipment rinsate and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

Nonconformances are summarized in Attachment A in Tables A-2 and A-3.

Sample results were qualified as follows:

For VOC samples:

| Blank type | Blank result | Sample result | Action for samples | |
|--|---------------------------|---|---|--|
| Method, Storage, Field, Trip, or Instrument* | Detects \leq LOQ | Not detected | No qualification | |
| | | < LOQ | Report sample LOQ value with a U | |
| | | \geq LOQ and \leq 2x LOQ | Report the sample result with a U** | |
| | > LOQ | \geq 2x the LOQ | No qualifications | |
| | | < LOQ | Report sample LOQ value with a U | |
| | | \geq LOQ and < blank contamination | Report the sample result with a U or reject the sample result as unusable R | |
| | | \geq LOQ and \geq blank contamination | If the result is \leq 2x blank result, report the sample result U.** If the result is $>$ 2x blank result, no qualification is required.** | |
| | | | | |
| | | | | |
| * Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L. | | | | |
| **Based on professional judgment. | | | | |

For TOC samples:

| Blank Type | Blank Result | Sample Result | Action for Samples |
|---------------------------|--------------------------|--|---|
| ICB/CCB (Positive) | \geq DL but \leq LOQ | Nondetect | No action |
| | | $>$ DL but \leq LOQ | Qualify as nondetect (U) at the LOQ |
| | | $>$ LOQ | Use professional judgment (see below [1]) |
| | >LOQ | $>$ DL but $<$ LOQ | Qualify as nondetect (U) at the LOQ |
| | | $>$ LOQ but $<$ ICB/CCB Result | Qualify at level of Blank Result with a "U" or Qualify result as unusable |
| | | $>$ ICB/CCB but $<$ 10x the ICB/CCB result | Qualify as estimated (J) |
| | | $>$ 10x ICB/CCB | No action is taken based on professional judgment |
| | | $>$ DL but \leq LOQ | Qualify as nondetect (U) at the LOQ |
| | | $>$ LOQ but $<$ 10x Blank Result | Qualify results as unusable |
| PB / EB/ FB (Positive) | > LOQ | $>$ 10x Blank Result | No action |
| | | Nondetect | No action |
| | | $>$ DL but \leq LOQ | Qualify as nondetect (U) at the LOQ |
| | | $>$ LOQ | Use professional judgment (see below [1]) |

[1] Establish an action level (AL) at 5x the blank contamination. If sample result is $<$ AL, qualify the reported result with a U.
LOQ - Limit of Quantitation.

Qualified sample results are shown in Table 1.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS/LCSD Results

The LCS/LCSD %Rs and/or relative percent recoveries (RPDs) were reviewed for conformance with the QC acceptance criteria.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|---|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) (LL = lower limit, UL = upper limit) | J | R |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than the reject sample results previously negated (U) on the basis of blank contamination.

Nonconformances are summarized in Attachment A in Table A-#. Qualified sample results are shown in Table 1.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-EB-090214 | WQ | TOTAL ORGANIC CARBON | | 1.0* | UG/L | UJ | bl |
| VPB154-EB-090214 | WQ | 2-BUTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-EB-090214 | WQ | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082914-838-840 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082914-838-840 | WG | ACETONE | 23 | 2.5 | UG/L | J | i |
| VPB154-GW-082914-838-840 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | be |
| VPB154-GW-082914-858-860 | WG | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-GW-082914-858-860 | WG | ACETONE | 22 | 2.5 | UG/L | J | i |
| VPB154-GW-082914-858-860 | WG | CARBON DISULFIDE | | 1.0* | UG/L | U | be |
| VPB154-GW-090214-908-910 | WG | 1,1,1-TRICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,1,2,2-TETRACHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,1,2-TRICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,1-DICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,1-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2,4-TRICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 38 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DIBROMOETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DICHLOROETHENE, TOTAL | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,2-DICHLOROPROPANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 1,3-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-090214-908-910 | WG | 1,4-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | 2-BUTANONE | | 120 | UG/L | UJ | mc,c |
| VPB154-GW-090214-908-910 | WG | 2-HEXANONE | | 120 | UG/L | UJ | mc,c |
| VPB154-GW-090214-908-910 | WG | 4-METHYL-2-PENTANONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | ACETONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | BENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | BROMODICHLOROMETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | BROMOFORM | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | BROMOMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CARBON DISULFIDE | | 50* | UG/L | UJ | mc,be |
| VPB154-GW-090214-908-910 | WG | CARBON TETRACHLORIDE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CHLOROETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CHLOROFORM | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CHLOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CIS-1,2-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CIS-1,3-DICHLOROPROPENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | CYCLOHEXANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | DIBROMOCHLOROMETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | DICHLORODIFLUOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | ETHYLBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | ISOPROPYLBENZENE | | 25 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|------|-------|-----------------------|-------------------|
| VPB154-GW-090214-908-910 | WG | M- AND P-XYLENE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | METHYL ACETATE | | 38 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | METHYL CYCLOHEXANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | METHYL TERT-BUTYL ETHER | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | METHYLENE CHLORIDE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | O-XYLENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | STYRENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TETRACHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TOLUENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TRANS-1,2-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TRANS-1,3-DICHLOROPROPENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TRICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | TRICHLOROFLUOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | VINYL CHLORIDE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090214-908-910 | WG | XYLENES, TOTAL | | 75 | UG/L | UJ | mc |
| VPB154-TRIP BLANK-090214 | WQ | 2-BUTANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-090214 | WQ | 2-HEXANONE | | 2.5 | UG/L | UJ | c |
| VPB154-TRIP BLANK-090214 | WQ | CARBON DISULFIDE | | 1.0* | UG/L | U | bl |

*LOQ

Attachment A**Nonconformance Summary Tables****Table A-1 -Continuing Calibration Verification Standard**

| CCV ID | Compound | % D | Limits |
|---|------------|-----|--------|
| WG149439-4 | 2-BUTANONE | -21 | ≤20% |
| | 2-HEXANONE | -24 | ≤20% |
| Associated samples: all samples in SDG SH7193 | | | |

Table A-2 - Lab Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------|----------------------|--------|------|-------|--------------------------|
| WG149729-1 | TOTAL ORGANIC CARBON | 0.50 | 0.50 | MG/L | VPB154-EB-090214 |
| WG149439-2 | CARBON DISULFIDE | 0.28 | 0.50 | UG/L | VPB154-TRIP BLANK-090214 |

Table A-3 - Field Blanks

| Blank ID | Compound | Result | LOD | Units | Associated Samples |
|------------------|------------------|--------|------|-------|--|
| VPB154-EB-090214 | CARBON DISULFIDE | 0.30 | 0.50 | UG/L | VPB154-GW-082914-838-840 VPB154-GW-082914-858-860 VPB154-GW-090214-908-910 |

Table A-4 - Lab Control Samples

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|----------|----------------|-------------|-------------|---|
| WG149439-1 | ACETONE | 146 | 40 | 140 | VPB154-GW-082914-838-840 VPB154-GW-082914-858-860 |

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|--------------------|--|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |
| mc | Method compliance nonconformance |



600 Technology Way
Scarborough, ME 04074
Tel: (207) 874-2400
Fax: (207) 775-4029

CHAIN of CUSTODY

PLEASE BEAR DOWN AND
PRINT LEGIBLY IN PEN

Page _____ of _____

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|--|--|--|------------------------------|------------------|--------------------------|------------------------------|----------------|--------------------------|------------------------------|----------------|
| Client <i>Revelation Consultants</i> | Contact <i>K. Viverette</i> | Phone # <i>(847) 457-4180 ()</i> | Fax # | | | | | | | |
| Address <i>160 Ryd Schoolhouse Rd.</i> | City <i>Chestnut Ridge</i> | State <i>NY</i> | Zip Code <i>10917</i> | | | | | | | |
| Purchase Order # | Proj. Name / No. <i>Nu IPP Netcharge / 60266528</i> | Katahdin Quote # | | | | | | | | |
| Bill (if different than above) | Address | | | | | | | | | |
| Sampler (Print / Sign) <i>Vince L. Viverette M.L.D.</i> | | | Copies To: | | | | | | | |
| LAB USE ONLY | WORK ORDER #: <i>SH 7193</i> | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | |
| KATAHDIN PROJECT NUMBER | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON |
| REMARKS: | | <i>UOC</i> | <i>TDC</i> | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | |
| TEMP°C <input type="checkbox"/> TEMP BLANK <input type="checkbox"/> INTACT <input type="checkbox"/> NOT INTACT | | | | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | | | | | | |
| | <i>VPD154-6L-082914-838870</i> | <i>8/29/14 / 1000</i> | <i>6W</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>VPB154-6L-082914-858-860</i> | <i>8/29/14 / 1000</i> | <i>6L</i> | <i>2</i> | <i>2</i> | | | | | |
| | <i>VPD154-EB-090214</i> | <i>9/2/14 / 1300</i> | <i>W</i> | <i>6</i> | <i>3</i> | <i>3</i> | | | | |
| | <i>VPD154-Trip Blank-090214</i> | <i>9/2/14 / 1430</i> | <i>W</i> | <i>3</i> | <i>3</i> | | | | | |
| | <i>VPB154-6L-090214-918-910</i> | <i>9/2/14 / 1520</i> | <i>6W</i> | <i>3</i> | <i>3</i> | | | | | |
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| COMMENTS <i>VPB154-6L-090214 - 9089-910 btl VPD154-6L-082914-858-860 - Very s. Hy.</i> | | | | | | | | | | |
| Relinquished By: (Signature) <i>[Signature]</i> | Date / Time <i>9/2/14 1700</i> | Received By: (Signature) <i>Fed EX</i> | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time |
| Relinquished By: (Signature) <i>[Signature]</i> | Date / Time <i>9/2/14 1700</i> | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time |

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN
SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-3
 Client ID: VPB154-EB-090214
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8843.D

Sample Date: 02-SEP-14
 Received Date: 03-SEP-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | J | 0.30 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

Page 1 of 2



ANALYTICAL SERVICES

Report of Analytical Results

Client: Rick Purdy
AECOM
701 Edgewater Drive
Wakefield, MA 01880

Lab Sample ID: SH7193-3

Report Date: 16-SEP-14

Client PO: 16518

Project: Navy Clean WE15-03-0

SDG: SH7193

Sample Description

VPB154-EB-090214

| Parameter | Result | Adj LOQ | Adj MDL | Adj LOD | Anal Method | QC Batch | Anal Date | Prep. Method | Prep. Date | Footnotes |
|----------------------|----------|---------|---------|---------|-------------|----------|--------------------|--------------|------------|-----------|
| Total Organic Carbon | 1047mg/L | 1.0 | 0.10 | .5 | SM5310B | WG149729 | 04-SEP-14 21:14:20 | N/A | N/A | |

1.0 0.5

✓ 1.0 ✓

Katahdin Analytical Services 0000135



Cert No EB7604

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-1
 Client ID: 154-082914-838-840
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8850.D

Sample Date: 29-AUG-14
 Received Date: 03-SEP-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- | 0.30 | 1.0 | ug/L | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- | 23 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | | 9.8 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | J | 0.33 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | +/- | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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12/24/14

Report of Analytical Results

Client: ENSAFE
Lab ID: SH7193-1
Client ID: 154-082914-838-840
Project: Navy Clean WE15-03-06 NW
SDG: SH7193
Lab File ID: C8850.D

Sample Date: 29-AUG-14 **Analysis Date:** 03-SEP-14
Received Date: 03-SEP-14 **Analyst:** REC
Extract Date: 03-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149439 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 96.1 | % | | | | | |
| Toluene-d8 | | 95.8 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 114. | % | | | | | |
| Dibromofluoromethane | | 100. | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-2
 Client ID: 154-082914-858-860
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8851.D

Sample Date: 29-AUG-14
 Received Date: 03-SEP-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- U | 0.48-1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | +/- J | 22 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | | 10 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | +/- U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | J | 0.67 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-2
 Client ID: 154-082914-858-860
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8851.D

Sample Date: 29-AUG-14 Analysis Date: 03-SEP-14
 Received Date: 03-SEP-14 Analyst: REC
 Extract Date: 03-SEP-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG149439 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 96.1 | % | | | | | |
| Toluene-d8 | | 95.3 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 114. | % | | | | | |
| Dibromofluoromethane | | 96.1 | % | | | | | |

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-5DL
 Client ID: 154-090214-908-910
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8844.D

Sample Date: 02-SEP-14
 Received Date: 03-SEP-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| Chloromethane | U | 50 | ug/L | 50 | 2 | 100 | 18. | 50. |
| Vinyl Chloride | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| Bromomethane | U | 50 | ug/L | 50 | 2 | 100 | 24. | 50. |
| Chloroethane | U | 50 | ug/L | 50 | 2 | 100 | 28. | 50. |
| Trichlorofluoromethane | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| 1,1-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| Carbon Disulfide | | 50 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Freon-113 | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Methylene Chloride | U | 120 | ug/L | 50 | 5 | 250 | 56. | 120 |
| Acetone | UL | 120 | ug/L | 50 | 5 | 250 | 110 | 120 |
| trans-1,2-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Methyl tert-butyl Ether | U | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| 1,1-Dichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| cis-1,2-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| Chloroform | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| 1,1,1-Trichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| 2-Butanone | U | 120 | ug/L | 50 | 5 | 250 | 66. | 120 |
| Cyclohexane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Carbon Tetrachloride | U | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| Benzene | U | 25 | ug/L | 50 | 1 | 50. | 13. | 25. |
| 1,2-Dichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| Trichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 14. | 25. |
| 1,2-Dichloropropane | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Bromodichloromethane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| cis-1,3-Dichloropropene | U | 25 | ug/L | 50 | 1 | 50. | 9.5 | 25. |
| Toluene | U | 25 | ug/L | 50 | 1 | 50. | 14. | 25. |
| 4-Methyl-2-Pentanone | U | 120 | ug/L | 50 | 5 | 250 | 66. | 120 |
| trans-1,3-Dichloropropene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| 1,1,2-Trichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Tetrachloroethene | U | 25 | ug/L | 50 | 1 | 50. | 20. | 25. |
| Dibromochloromethane | U | 25 | ug/L | 50 | 1 | 50. | 15. | 25. |
| 2-Hexanone | U | 120 | ug/L | 50 | 5 | 250 | 85. | 120 |
| Chlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| Ethylbenzene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7193-5DL
Client ID: 154-090214-908-910
Project: Navy Clean WE15-03-06 NW
SDG: SH7193
Lab File ID: C8844.D

Sample Date: 02-SEP-14 **Analysis Date:** 03-SEP-14
Received Date: 03-SEP-14 **Analyst:** REC
Extract Date: 03-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149439 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | ✓ | 75 | ug/L | 50 | 3 | 150 | 12. | 75. |
| Styrene | ✓ | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Bromoform | ✓ | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Isopropylbenzene | ✓ | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| 1,1,2,2-Tetrachloroethane | ✓ | 25 | ug/L | 50 | 1 | 50. | 19. | 25. |
| 1,3-Dichlorobenzene | ✓ | 25 | ug/L | 50 | 1 | 50. | 13. | 25. |
| 1,4-Dichlorobenzene | ✓ | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| 1,2-Dichlorobenzene | ✓ | 25 | ug/L | 50 | 1 | 50. | 7.5 | 25. |
| 1,2,4-Trichlorobenzene | ✓ | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| Methyl Acetate | ✓ | 38 | ug/L | 50 | 1 | 50. | 26. | 38. |
| Methylcyclohexane | ✓ | 25 | ug/L | 50 | 1 | 50. | 15. | 25. |
| o-Xylene | ✓ | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| M+P-Xylenes | ✓ | 50 | ug/L | 50 | 2 | 100 | 30. | 50. |
| 1,2-Dichloroethylene (Total) | ✓ | 50 | ug/L | 50 | 2 | 100 | 10. | 50. |
| 1,2-Dibromoethane | ✓ | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| 1,2-Dibromo-3-Chloropropane | ✓ | 38 | ug/L | 50 | 1 | 50. | 25. | 38. |
| P-Bromofluorobenzene | | 94.2 | % | | | | | |
| Toluene-d8 | | 94.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 106. | % | | | | | |
| Dibromofluoromethane | | 96.4 | % | | | | | |

Q.2/22/14

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7193-4
 Client ID: VPB154-TB-090214
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7193
 Lab File ID: C8842.D

Sample Date: 02-SEP-14
 Received Date: 03-SEP-14
 Extract Date: 03-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149439

Analysis Date: 03-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|----------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | +/- U | 0.27-1.0 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | UL | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | +/- U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | +/- U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7193-4
Client ID: VPB154-TB-090214
Project: Navy Clean WE15-03-06 NW
SDG: SH7193
Lab File ID: C8842.D

Sample Date: 02-SEP-14 **Analysis Date:** 03-SEP-14
Received Date: 03-SEP-14 **Analyst:** REC
Extract Date: 03-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149439 **Report Date:** 04-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 90.4 | % | | | | | |
| Toluene-d8 | | 91.5 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 103. | % | | | | | |
| Dibromofluoromethane | | 93.2 | % | | | | | |



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Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage

Laboratory: Katahdin Analytical

Service Request: SH7248

Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS)

Validation Level: 3

AECOM Project 60266526.SA.DV

Number:

Prepared by: Dawn Brule/RESCON Completed on: 12/05/2014

Reviewed by: Lori Herberich/RESCON File Name: SH7248_8260B

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on September 3, 2014.

| Sample ID | Matrix/Sample Type |
|--------------------------|--------------------|
| VPB154-GW-090314-918-920 | Groundwater |
| VPB154-TRIP BLANK-090314 | Trip Blank |

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- Data completeness (chain-of-custody [COC])/sample integrity
- Holding times and sample preservation
- GC/MS performance checks
- Initial calibration/continuing calibration verification
- Laboratory blanks/equipment blanks/trip blanks
- Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- X Laboratory control sample (LCS) results
- NA Field duplicate results
- Internal standard results
- Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (✗) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Selected data points were estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

The vials of sample VPB154-GW-090314-918-920 were mostly soil and had very little standing water. Therefore, each vial was decanted and composited into one vial and analyzed. Due to limited sample volume, the sample was analyzed at a 1:50 dilution. Positive and nondetect results for this sample were qualified as estimated (J and UJ, respectively), due to possible loss of sample integrity during the decanting procedure. Qualified sample results are shown in Table 1.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r^2), and/or response factor method acceptance criteria were met;

- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

The QC acceptance criteria were met.

Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip blanks were evaluated as to whether there were contaminants detected above the detection limit (DL).

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method, equipment rinsate and trip blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

The QC acceptance criteria were met and/or qualification of the sample results was not required.

Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

MS/MSD Results

MS/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria. Nonconformances are summarized in Attachment A in Table A-1.

Data qualification to the analytes associated with the specific LCS %Rs or RPDs was as follows:

| Nonconformances ¹ | Action | |
|---|--------------------|-----------------------|
| | Detected Compounds | Nondetected Compounds |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% (see note 1) (LL = lower limit, UL = upper limit) | J | R |

Notes:

1. Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather than reject sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1.



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Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Internal Standard Results

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-090314-918-920 | WG | 1,1,1-TRICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,1,2,2-TETRACHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,1,2-TRICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,1-DICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,1-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2,4-TRICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2-DIBROMO-3-CHLOROPROPANE | | 38 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2-DIBROMOETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2-DICHLOROETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,2-DICHLOROETHENE, TOTAL | | 50 | UG/L | UJ | mc,l |
| VPB154-GW-090314-918-920 | WG | 1,2-DICHLOROPROPANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,3-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 1,4-DICHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 2-BUTANONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 2-HEXANONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | 4-METHYL-2-PENTANONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | ACETONE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | BENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | BROMODICHLOROMETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | BROMOFORM | | 25 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|-------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-090314-918-920 | WG | BROMOMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CARBON DISULFIDE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CARBON TETRACHLORIDE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CHLOROBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CHLOROETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CHLOROFORM | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CHLOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CIS-1,2-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CIS-1,3-DICHLOROPROPENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | CYCLOHEXANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | DIBROMOCHLOROMETHANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | DICHLORODIFLUOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | ETHYLBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | ISOPROPYLBENZENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | M- AND P-XYLENE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | METHYL ACETATE | | 38 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | METHYL CYCLOHEXANE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | METHYL TERT-BUTYL ETHER | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | METHYLENE CHLORIDE | | 120 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | O-XYLENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | STYRENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | TETRACHLOROETHENE | | 25 | UG/L | UJ | mc |

| Sample ID | Matrix | Compound | Result | LOD | Units | Validation Qualifiers | Validation Reason |
|--------------------------|--------|---------------------------|--------|-----|-------|-----------------------|-------------------|
| VPB154-GW-090314-918-920 | WG | TOLUENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | TRANS-1,2-DICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | TRANS-1,3-DICHLOROPROPENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | TRICHLOROETHENE | | 25 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | TRICHLOROFLUOROMETHANE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | VINYL CHLORIDE | | 50 | UG/L | UJ | mc |
| VPB154-GW-090314-918-920 | WG | XYLENES, TOTAL | | 75 | UG/L | UJ | mc |
| VPB154-TRIP BLANK-090314 | WQ | 1,2-DICHLOROETHENE, TOTAL | | 1.0 | UG/L | UJ | I |

Attachment A**Nonconformance Summary Tables****Table A-1 - Lab Control Samples**

| LCS ID | Compound | LCS % Recovery | Lower Limit | Upper Limit | Associated Samples |
|------------|---------------------------|----------------|-------------|-------------|--|
| WG149482-1 | 1,2-DICHLOROETHENE, TOTAL | 80.7 | 84 | 121 | VPB154-GW-090314-918-920 VPB154-TRIP BLANK-090314 |

Attachment B**Qualifier Codes and Explanations**

| Qualifier | Explanation |
|-----------|---|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified. |

Attachment C**Reason Codes and Explanations**

| Reason Code | Explanation |
|-------------|--|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| c | Calibration issue |
| co | Analyte carryover |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate RPDs |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration (EMPC) |
| l | LCS or OPR recoveries |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate RPDs |
| lp | Laboratory control sample/laboratory control sample duplicate RPDs |
| m | Matrix spike recovery |
| md | Matrix spike/matrix spike duplicate RPDs |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column RPD |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | ICS results |
| mc | Method compliance nonconformance |



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Page 1 of 1

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| Client Resolution Consultant | Contact E. Vivera | Phone # (845)425-4180 | Fax # | | | | | | | | | | | |
| Address 106 Red Schoolhouse Rd | City Chestnut Ridge | State NY | Zip Code 10977 | | | | | | | | | | | |
| Purchase Order # | Proj. Name / No. NL1RP Bethpage / 60266526 | Katahdin Quote # | | | | | | | | | | | | |
| Bill (if different than above) | Address | | | | | | | | | | | | | |
| Sampler (Print / Sign) <i>Vincent Vassichio / V.V.</i> | | | Copies To: | | | | | | | | | | | |
| LAB USE ONLY | WORK ORDER #: SH7248 | ANALYSIS AND CONTAINER TYPE PRESERVATIVES | | | | | | | | | | | | |
| KATAHDIN PROJECT NUMBER | | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | Filt. OY ON | | | | |
| REMARKS: | | <i>VOCs</i> | | | | | | | | | | | | |
| SHIPPING INFO: <input type="checkbox"/> FED EX <input type="checkbox"/> UPS <input type="checkbox"/> CLIENT | | | | | | | | | | | | | | |
| AIRBILL NO: | | | | | | | | | | | | | | |
| TEMP'C | | <input type="checkbox"/> TEMP BLANK | <input type="checkbox"/> INTACT | <input type="checkbox"/> NOT INTACT | | | | | | | | | | |
| * | Sample Description | Date / Time coll'd | Matrix | No. of Cntrs. | | | | | | | | | | |
| | UPD154-6W-090314-920 | 9/3/14 605 | 6W | 3 | | | | | | | | | | |
| | UPD154-TripN61c-090314 | 9/3/14 / 1430 | w | 3 | | | | | | | | | | |
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| COMMENTS <i>Sample Very Silty</i> | | | | | | | | | | | | | | |
| Relinquished By: (Signature) <i>V.V.</i> | Date / Time 9/3/14 / 1600 | Received By: (Signature) Fed EX | Relinquished By: (Signature) | Date / Time | Received By: (Signature) <i>7/4/14 0915</i> | | | | | | | | | |
| Relinquished By: (Signature) | Date / Time | Received By: (Signature) | Relinquished By: (Signature) | Date / Time | Received By: (Signature) | | | | | | | | | |

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KAS-COC1

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ED 002631A 00004611-00280

Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7248-1DL
 Client ID: 154-090314-918-920
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7248
 Lab File ID: C8867.D

Sample Date: 03-SEP-14
 Received Date: 04-SEP-14
 Extract Date: 04-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149482

Analysis Date: 04-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 05-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| Chloromethane | U | 50 | ug/L | 50 | 2 | 100 | 18. | 50. |
| Vinyl Chloride | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| Bromomethane | U | 50 | ug/L | 50 | 2 | 100 | 24. | 50. |
| Chloroethane | U | 50 | ug/L | 50 | 2 | 100 | 28. | 50. |
| Trichlorofluoromethane | U | 50 | ug/L | 50 | 2 | 100 | 12. | 50. |
| 1,1-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| Carbon Disulfide | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Freon-113 | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Methylene Chloride | U | 120 | ug/L | 50 | 5 | 250 | 56. | 120 |
| Acetone | UL | 120 | ug/L | 50 | 5 | 250 | 110 | 120 |
| trans-1,2-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Methyl tert-butyl Ether | U | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| 1,1-Dichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| cis-1,2-Dichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| Chloroform | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| 1,1,1-Trichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| 2-Butanone | U | 120 | ug/L | 50 | 5 | 250 | 66. | 120 |
| Cyclohexane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Carbon Tetrachloride | U | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| Benzene | U | 25 | ug/L | 50 | 1 | 50. | 13. | 25. |
| 1,2-Dichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| Trichloroethene | U | 25 | ug/L | 50 | 1 | 50. | 14. | 25. |
| 1,2-Dichloropropane | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Bromodichloromethane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| cis-1,3-Dichloropropene | U | 25 | ug/L | 50 | 1 | 50. | 9.5 | 25. |
| Toluene | U | 25 | ug/L | 50 | 1 | 50. | 14. | 25. |
| 4-Methyl-2-Pentanone | U | 120 | ug/L | 50 | 5 | 250 | 66. | 120 |
| trans-1,3-Dichloropropene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |
| 1,1,2-Trichloroethane | U | 25 | ug/L | 50 | 1 | 50. | 16. | 25. |
| Tetrachloroethene | U | 25 | ug/L | 50 | 1 | 50. | 20. | 25. |
| Dibromochloromethane | U | 25 | ug/L | 50 | 1 | 50. | 15. | 25. |
| 2-Hexanone | U | 120 | ug/L | 50 | 5 | 250 | 85. | 120 |
| Chlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| Ethylbenzene | U | 25 | ug/L | 50 | 1 | 50. | 10. | 25. |

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Report of Analytical Results

Client: ENSAFE
 Lab ID: SH7248-1DL
 Client ID: 154-090314-918-920
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7248
 Lab File ID: C8867.D

Sample Date: 03-SEP-14 Analysis Date: 04-SEP-14
 Received Date: 04-SEP-14 Analyst: REC
 Extract Date: 04-SEP-14 Analysis Method: SW846 8260C
 Extracted By: REC Matrix: AQ
 Extraction Method: SW846 5030 % Solids: NA
 Lab Prep Batch: WG149482 Report Date: 05-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 75 | ug/L | 50 | 3 | 150 | 12. | 75. |
| Styrene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Bromoform | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| Isopropylbenzene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| 1,1,2,2-Tetrachloroethane | U | 25 | ug/L | 50 | 1 | 50. | 19. | 25. |
| 1,3-Dichlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 13. | 25. |
| 1,4-Dichlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| 1,2-Dichlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 7.5 | 25. |
| 1,2,4-Trichlorobenzene | U | 25 | ug/L | 50 | 1 | 50. | 18. | 25. |
| Methyl Acetate | U | 38 | ug/L | 50 | 1 | 50. | 26. | 38. |
| Methylcyclohexane | U | 25 | ug/L | 50 | 1 | 50. | 15. | 25. |
| o-Xylene | U | 25 | ug/L | 50 | 1 | 50. | 12. | 25. |
| M+P-Xylenes | U | 50 | ug/L | 50 | 2 | 100 | 30. | 50. |
| 1,2-Dichloroethylene (Total) | UL | 50 | ug/L | 50 | 2 | 100 | 10. | 50. |
| 1,2-Dibromoethane | U | 25 | ug/L | 50 | 1 | 50. | 11. | 25. |
| 1,2-Dibromo-3-Chloropropane | U | 38 | ug/L | 50 | 1 | 50. | 25. | 38. |
| P-Bromofluorobenzene | | 91.0 | % | | | | | |
| Toluene-d8 | | 91.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 105. | % | | | | | |
| Dibromofluoromethane | | 90.2 | % | | | | | |



Report of Analytical Results

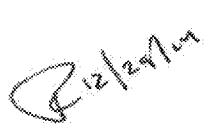
Client: ENSAFE
 Lab ID: SH7248-2
 Client ID: VPB154-TB-090314
 Project: Navy Clean WE15-03-06 NW
 SDG: SH7248
 Lab File ID: C8864.D

Sample Date: 03-SEP-14
 Received Date: 04-SEP-14
 Extract Date: 04-SEP-14
 Extracted By: REC
 Extraction Method: SW846 5030
 Lab Prep Batch: WG149482

Analysis Date: 04-SEP-14
 Analyst: REC
 Analysis Method: SW846 8260C
 Matrix: AQ
 % Solids: NA
 Report Date: 05-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|---------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Dichlorodifluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| Chloromethane | J | 0.55 | ug/L | 1 | 2 | 2.0 | 0.36 | 1.0 |
| Vinyl Chloride | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.25 | 1.0 |
| Bromomethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.49 | 1.0 |
| Chloroethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.55 | 1.0 |
| Trichlorofluoromethane | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.24 | 1.0 |
| 1,1-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.35 | 0.50 |
| Carbon Disulfide | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Freon-113 | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Methylene Chloride | J | 1.2 | ug/L | 1 | 5 | 5.0 | 1.1 | 2.5 |
| Acetone | DL | 2.5 | ug/L | 1 | 5 | 5.0 | 2.2 | 2.5 |
| trans-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Methyl tert-butyl Ether | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.36 | 0.50 |
| 1,1-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| cis-1,2-Dichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |
| Chloroform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.32 | 0.50 |
| 1,1,1-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 2-Butanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| Cyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.31 | 0.50 |
| Carbon Tetrachloride | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Benzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,2-Dichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| Trichloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.28 | 0.50 |
| 1,2-Dichloropropane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| Bromodichloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| cis-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.19 | 0.50 |
| Toluene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.27 | 0.50 |
| 4-Methyl-2-Pentanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.3 | 2.5 |
| trans-1,3-Dichloropropene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.20 | 0.50 |
| 1,1,2-Trichloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.33 | 0.50 |
| Tetrachloroethene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.40 | 0.50 |
| Dibromochloromethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| 2-Hexanone | U | 2.5 | ug/L | 1 | 5 | 5.0 | 1.7 | 2.5 |
| Chlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| Ethylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.21 | 0.50 |

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Report of Analytical Results

Client: ENSAFE
Lab ID: SH7248-2
Client ID: VPB154-TB-090314
Project: Navy Clean WE15-03-06 NW
SDG: SH7248
Lab File ID: C8864.D

Sample Date: 03-SEP-14 **Analysis Date:** 04-SEP-14
Received Date: 04-SEP-14 **Analyst:** REC
Extract Date: 04-SEP-14 **Analysis Method:** SW846 8260C
Extracted By: REC **Matrix:** AQ
Extraction Method: SW846 5030 **% Solids:** NA
Lab Prep Batch: WG149482 **Report Date:** 05-SEP-14

| Compound | Qualifier | Result | Units | Dilution | LOQ | ADJ LOQ | ADJ MDL | ADJ LOD |
|------------------------------|-----------|--------|-------|----------|-----|---------|---------|---------|
| Xylenes (total) | U | 1.5 | ug/L | 1 | 3 | 3.0 | 0.25 | 1.5 |
| Styrene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Bromoform | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| Isopropylbenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.23 | 0.50 |
| 1,1,2,2-Tetrachloroethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.38 | 0.50 |
| 1,3-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.26 | 0.50 |
| 1,4-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.24 | 0.50 |
| 1,2-Dichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.15 | 0.50 |
| 1,2,4-Trichlorobenzene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.37 | 0.50 |
| Methyl Acetate | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.53 | 0.75 |
| Methylcyclohexane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.30 | 0.50 |
| o-Xylene | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.25 | 0.50 |
| M+P-Xylenes | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.59 | 1.0 |
| 1,2-Dichloroethylene (Total) | U | 1.0 | ug/L | 1 | 2 | 2.0 | 0.21 | 1.0 |
| 1,2-Dibromoethane | U | 0.50 | ug/L | 1 | 1 | 1.0 | 0.22 | 0.50 |
| 1,2-Dibromo-3-Chloropropane | U | 0.75 | ug/L | 1 | 1 | 1.0 | 0.50 | 0.75 |
| P-Bromofluorobenzene | | 93.6 | % | | | | | |
| Toluene-d8 | | 92.2 | % | | | | | |
| 1,2-Dichloroethane-d4 | | 109. | % | | | | | |
| Dibromofluoromethane | | 93.7 | % | | | | | |



Section 5
VPB 154 Analytical Data Table

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|----------------------------|-----------------------------|------------------------------|--------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/7/2014 | 8/7/2014 | 8/7/2014 | 8/7/2014 |
| Sample ID | | VPB154-GW-080714- 58-60 | VPB154-GW-080714- 98-100 | VPB154-GW-080714- 148-150 | VPB154-GWD-080714 |
| Sample Interval | (Note 1) | 58 - 60 ft | 98 - 100 ft | 148 - 150 ft | 148 - 150 ft |
| Sample type code | | N | N | N | FD |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | 2.9 J |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | 7.7 | 8.3 | 12 | 13 |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CARBON DISULFIDE | 60 | < 0.50 U | < 0.50 U | < 0.50 U | < 1.0 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | < 0.50 U | < 0.50 U | 0.33 J | < 0.50 U |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TOLUENE | 5 | 0.30 J | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROFUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/8/2014 | 8/8/2014 | 8/11/2014 | 8/12/2014 |
| Sample ID | | VPB154-GW-080814- 198-200 | VPB154-GW-080814- 228-230 | VPB154-GW-081114- 238-240 | VPB154-GW-081214- 258-260 |
| Sample Interval | (Note 1) | 198 - 200 ft | 228 - 230 ft | 238 - 240 ft | 258 - 260 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | 0.21 J | < 0.50 U |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 1.7 | 10 | 14 | < 0.50 U |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 0.50 U | 0.61 J | 0.81 J | < 0.50 U |
| 1,1-DICHLOROETHENE | 5 | < 0.50 U | 1.2 J | 1.5 J | < 0.50 U |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 0.89 J | 2.6 | 4.1 | < 1.0 U |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | 20 J | 2.7 J | 3.4 J | 7.0 J |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 UJ |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 UJ |
| ACETONE | 50 | 72 J | 9.8 J | 16 J | 48 J |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CARBON DISULFIDE | 60 | 0.26 J | 0.25 J | 0.30 J | < 1.0 UJ |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | < 0.50 U | 1.4 | 0.86 J | < 0.50 U |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | 0.89 J | 2.6 | 4.1 | < 0.50 U |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | < 1.0 U | 0.70 J | 0.35 J | < 1.0 U |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | 1.0 | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 3.2 | 4.8 | 5.1 | < 0.50 U |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 31 | 180 | 190 | < 0.50 U |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/12/2014 | 8/12/2014 | 8/12/2014 | 8/14/2014 |
| Sample ID | | VPB154-GW-081214- 278-280 | VPB154-GW-081214- 298-300 | VPB154-GW-081214- 318-320 | VPB154-GW-081414- 338-340 |
| Sample Interval | (Note 1) | 278 - 280 ft | 298 - 300 ft | 318 - 320 ft | 338 - 340 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 16 | 30 | 16 J | 13 |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 1.0 U | < 1.2 U | < 1.0 U | < 1.0 U |
| 1,1-DICHLOROETHENE | 5 | 2.2 J | 3.3 J | 2.2 J | 1.5 J |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | 0.34 J |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 4.8 | 5.2 | 2.8 J | 2.3 |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | 3.7 J | < 2.5 U | 1.9 J | 1.5 J |
| 2-HEXANONE | 50 | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ | < 2.5 U |
| ACETONE | 50 | < 15 UJ | < 5.0 UJ | < 5.0 UJ | < 5.0 UJ |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CARBON DISULFIDE | 60 | < 1.0 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | 4.8 | 5.2 | 2.8 J | 2.3 |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | 0.33 J | 0.74 J | 0.79 J | 1.3 J |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYL TERT-BUTYL ETHER | 10 | 0.45 J | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 7.1 | 8.0 | 7.5 J | < 0.50 U |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 200 | 200 | 180 | 26 |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/14/2014 | 8/14/2014 | 8/15/2014 | 8/15/2014 |
| Sample ID | | VPB154-GW-081414- 358-360 | VPB154-GW-081414- 378-380 | VPB154-GW-081514- 398-400 | VPB154-GW-081514- 418-420 |
| Sample Interval | (Note 1) | 358 - 360 ft | 378 - 380 ft | 398 - 400 ft | 418 - 420 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | 0.25 J |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | < 0.50 UJ | 0.41 J | < 0.50 UJ | 16 J |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | 0.56 J |
| 1,1-DICHLOROETHENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | 1.9 |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 UJ | < 0.75 U | < 0.75 U | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | 2.6 |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | 3.8 J | < 2.5 U | 2.7 J | < 2.5 U |
| 2-HEXANONE | 50 | < 2.5 UJ | < 2.5 U | < 2.5 U | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 UJ | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | < 17 UJ | < 5.0 UJ | 12 | 8.3 |
| BENZENE | 1 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| CARBON DISULFIDE | 60 | < 1.0 UJ | < 1.0 UJ | < 1.0 U | < 1.0 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | < 0.50 UJ | < 0.50 U | < 0.50 U | 0.87 J |
| CHLOROMETHANE | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | 2.6 |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | 1.0 J |
| ETHYLBENZENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 UJ | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 UJ | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | 4.8 |
| TOLUENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 UJ | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 1.5 J | 0.34 J | < 0.50 U | 170 |
| TRICLOROFUOROMETHANE | 5 | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 UJ | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 UJ | < 1.5 U | < 1.5 U | < 1.5 U |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/15/2014 | 8/18/2014 | 8/18/2014 | 8/19/2014 |
| Sample ID | | VPB154-GW-081514- 438-440 | VPB154-GW-081814- 458-460 | VPB154-GW-081814- 483-485 | VPB154-GW-081914- 503-505 |
| Sample Interval | (Note 1) | 438 - 440 ft | 458 - 460 ft | 483 - 485 ft | 503 - 505 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | 0.26 J | < 0.50 U | < 0.50 U | 0.53 J |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 14 J | 10 J | 7.7 J | 20 J |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | 0.40 J |
| 1,1-DICHLOROETHANE | 5 | 0.49 J | 0.44 J | 0.24 J | < 0.50 U |
| 1,1-DICHLOROETHENE | 5 | 1.7 | 1.1 | 0.71 J | 2.9 J |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 UJ |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 3.0 | 2.4 | 1.6 J | 2.2 |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | 4.9 J | 4.0 J | < 2.5 U | < 2.5 U |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | 17 | 22 | 9.7 | 8.4 J |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 UJ |
| CARBON DISULFIDE | 60 | < 1.0 U | < 1.0 U | < 0.50 U | < 0.50 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | 0.76 J | 0.43 J | < 0.50 U | 1.4 |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 2.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | 3.0 | 2.4 | 1.6 | 2.2 |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | 0.97 J | 0.43 J | 0.41 J | 2.0 |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 5.0 | 5.4 | 6.6 | < 0.50 U |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 190 | 180 | 160 | 160 |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | 0.28 J | 0.34 J | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/19/2014 | 8/20/2014 | 8/20/2014 | 8/20/2014 |
| Sample ID | | VPB154-GW-081914- 518-520 | VPB154-GW-082014- 538-540 | VPB154-GW-082014- 558-560 | VPB154-GW-082014- 578-580 |
| Sample Interval | (Note 1) | 518 - 520 ft | 538 - 540 ft | 558 - 560 ft | 578 - 580 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | 0.41 J | 0.87 J | 1.1 | 2.6 |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 26 J | 61 J | 57 J | 120 J |
| 1,1,2-TRICHLOROETHANE | 1 | 0.50 J | 1.1 | 0.98 J | 1.8 |
| 1,1-DICHLOROETHANE | 5 | 0.79 J | 2.6 | 2.6 | 3.6 |
| 1,1-DICHLOROETHENE | 5 | 5.5 J | 18 J | 19 J | 39 J |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 2.3 | 4.4 | 3.9 | 5.2 |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | < 2.5 U | 3.8 J | 2.4 J | < 2.5 U |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | 8.3 J | 20 J | 9.6 J | 8.3 J |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ |
| CARBON DISULFIDE | 60 | < 1.0 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | 0.74 J | 1.1 | 0.98 J | 1.2 |
| CHLOROMETHANE | 5 | < 1.0 U | < 2.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | 2.3 | 4.4 | 3.9 | 5.2 |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | 0.66 J | 0.49 J | 0.49 J | 0.79 J |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 3.4 | 12 | 10 | 8.4 |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 340 | 970 | 900 | 1800 J |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | 0.42 J | 0.47 J | 0.80 J |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|---------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/21/2014 | 8/21/2014 | 8/21/2014 | 8/21/2014 |
| Sample ID | | VPB154-GW-082114- 598-600 | VPB154-GW-082114- 618-620 | VPB154-GW-082114- 638-640 | VPB154-GWD-082114 |
| Sample Interval | (Note 1) | 598 - 600 ft | 618 - 620 ft | 638 - 640 ft | 638 - 640 ft |
| Sample type code | | N | N | N | FD |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | 1.3 | 3.7 | 2.5 | 2.5 |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 41 J | 70 J | 69 J | 70 J |
| 1,1,2-TRICHLOROETHANE | 1 | 1.1 | 1.6 | 1.6 | 1.6 |
| 1,1-DICHLOROETHANE | 5 | 1.8 | 4.7 | 4.5 | 4.4 |
| 1,1-DICHLOROETHENE | 5 | 14 J | 44 J | 29 J | 28 J |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ | < 0.75 UJ |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 2.2 | 3.1 | 4.4 | 4.1 |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| 2-BUTANONE | 50 | 4.5 J | 2.3 J | 1.4 J | 1.5 J |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | 12 J | 10 J | 6.2 J | 6.2 J |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ |
| CARBON DISULFIDE | 60 | < 1.0 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CHLOROFORM | 7 | 1.6 | 1.5 | 1.3 | 1.3 |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | 2.2 | 3.1 | 4.4 | 4.1 |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | 0.87 J | 0.54 J | 0.39 J | 0.34 J |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 0.40 J | 1.7 | 5.6 | 5.6 |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 350 J | 810 J | 1800 J | 1600 J |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | 0.37 J | 0.40 J | 0.42 J |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 1.5 U |

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/22/2014 | 8/22/2014 | 8/25/2014 | 8/25/2014 |
| Sample ID | | VPB154-GW-082214- 663-665 | VPB154-GW-082214- 678-680 | VPB154-GW-082514- 698-700 | VPB154-GW-082514- 718-720 |
| Sample Interval | (Note 1) | 663 - 665 ft | 678 - 680 ft | 698 - 700 ft | 718 - 720 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | 2.0 J | 0.37 J | 0.60 J | < 10 UJ |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 84 J | 21 J | 30 J | < 10 UJ |
| 1,1,2-TRICHLOROETHANE | 1 | 1.2 J | 0.53 J | 0.68 J | < 10 UJ |
| 1,1-DICHLOROETHANE | 5 | 2.6 J | 0.97 J | 1.5 J | < 10 UJ |
| 1,1-DICHLOROETHENE | 5 | 16 J | 3.2 J | 6.0 J | < 10 UJ |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 UJ | < 15 UJ |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 3.9 J | 2.1 | 2.9 J | < 20 UJ |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| 2-BUTANONE | 50 | 2.4 J | < 2.5 U | 5.3 J | < 50 UJ |
| 2-HEXANONE | 50 | < 2.5 U | < 2.5 U | < 2.5 UJ | < 50 UJ |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 UJ | < 50 UJ |
| ACETONE | 50 | 10 J | 8.7 J | 26 J | < 50 UJ |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| CARBON DISULFIDE | 60 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ | < 10 UJ |
| CARBON TETRACHLORIDE | 5 | 2.4 J | < 0.50 U | < 0.50 UJ | < 10 UJ |
| CHLOROBENZENE | 5 | < 0.50 U | 0.39 J | < 0.50 UJ | < 10 UJ |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| CHLOROFORM | 7 | 1.6 J | 0.78 J | 1.2 J | < 10 UJ |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| CIS-1,2-DICHLOROETHENE | 5 | 3.9 J | 2.1 | 2.9 J | < 10 UJ |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| DICHLORODIFLUOROMETHANE | 5 | 0.26 J | < 1.0 U | < 1.0 UJ | < 20 UJ |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 UJ | < 15 UJ |
| METHYL CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ | < 10 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 UJ | < 50 UJ |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| TETRACHLOROETHENE | 5 | < 0.50 U | 2.8 | 4.6 J | < 10 UJ |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 UJ | < 10 UJ |
| TRICHLOROETHENE | 5 | 600 | 550 | 700 J | < 10 UJ |
| TRICLOROFUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 UJ | < 20 UJ |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 UJ | < 30 UJ |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/25/2014 | 8/27/2014 | 8/27/2014 | 8/28/2014 |
| Sample ID | | VPB154-GW-082514- 738-740 | VPB154-GW-082714- 763-765 | VPB154-GW-082714- 778-780 | VPB154-GW-082814- 798-800 |
| Sample Interval | (Note 1) | 738 - 740 ft | 763 - 765 ft | 778 - 780 ft | 798 - 800 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,1,2-TRICHLOROETHANE | 1 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,1-DICHLOROETHENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,2,4-TRICHLOROBENZENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 1.5 UJ | < 0.75 U | < 1.5 UJ | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| 1,2-DICHLOROPROPANE | 1 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| 2-BUTANONE | 50 | < 5.0 UJ | < 2.5 U | < 5.0 UJ | < 2.5 UJ |
| 2-HEXANONE | 50 | < 5.0 UJ | < 2.5 U | < 5.0 UJ | < 2.5 UJ |
| 4-METHYL-2-PENTANONE | NL | < 5.0 UJ | < 2.5 U | < 5.0 UJ | < 2.5 U |
| ACETONE | 50 | 15 J | 13 J | 38 J | 13 J |
| BENZENE | 1 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| BROMOFORM | 50 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| BROMOMETHANE | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| CARBON DISULFIDE | 60 | < 1.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| CARBON TETRACHLORIDE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| CHLOROBENZENE | 5 | 0.69 J | < 0.50 U | < 1.0 UJ | < 0.50 U |
| CHLOROETHANE | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| CHLOROFORM | 7 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| CHLOROMETHANE | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| CIS-1,2-DICHLOROETHENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| CYCLOHEXANE | NL | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| DIBROMOCHLOROMETHANE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| ETHYLBENZENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| M- AND P-XYLENE | NL | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| METHYL ACETATE | NL | < 1.5 UJ | < 0.75 U | < 1.5 UJ | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| METHYL TERT-BUTYL ETHER | 10 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 5.0 UJ | < 2.5 U | < 5.0 UJ | < 2.5 U |
| O-XYLENE | NL | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| STYRENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TETRACHLOROETHENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TOLUENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TRANS-1,2-DICHLOROETHENE | 5 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 1.0 UJ | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TRICHLOROETHENE | 5 | 1.4 J | < 0.50 U | < 1.0 UJ | < 0.50 U |
| TRICLOROFUOROMETHANE | 5 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| VINYL CHLORIDE | 2 | < 2.0 UJ | < 1.0 U | < 2.0 UJ | < 1.0 U |
| XYLENES, TOTAL | 5 | < 3.0 UJ | < 1.5 U | < 3.0 UJ | < 1.5 U |

Vertical Profile Boring 154 Analytical Data Table
 Naval Weapons Industrial Reserve Plant
 Bethpage - Bethpage, New York

Resolution Consultants

| Location | | VPB154 | VPB154 | VPB154 | VPB154 |
|---------------------------------------|--|------------------------------|------------------------------|------------------------------|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value | 8/28/2014 | 8/29/2014 | 8/29/2014 | 9/2/2014 |
| Sample ID | | VPB154-GW-082814- 818-820 | VPB154-GW-082914- 838-840 | VPB154-GW-082914- 858-860 | VPB154-GW-090214- 908-910 |
| Sample Interval | (Note 1) | 818 - 820 ft | 838 - 840 ft | 858 - 860 ft | 908 - 910 ft |
| Sample type code | | N | N | N | N |
| VOC 8260C (ug/L) | | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,1,2-TRICHLOROETHANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,1-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,1-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U | < 38 UJ |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,2-DICHLOROETHENE, TOTAL | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| 2-BUTANONE | 50 | < 2.5 UJ | 9.8 | 10 | < 120 UJ |
| 2-HEXANONE | 50 | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ | < 120 UJ |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U | < 120 UJ |
| ACETONE | 50 | 5.9 J | 23 J | 22 J | < 120 UJ |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| BROMOMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| CARBON DISULFIDE | 60 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | 0.67 J | < 25 UJ |
| CHLOROETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| CHLOROFORM | 7 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| CHLOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| CIS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| DICHLORODIFLUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U | < 38 UJ |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U | < 120 UJ |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| TETRACHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| TOLUENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U | < 25 UJ |
| TRICHLOROETHENE | 5 | < 0.50 U | 0.33 J | < 0.50 U | < 25 UJ |
| TRICHLOROFUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U | < 50 UJ |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U | < 75 UJ |

| Location | | VPB154 |
|---------------------------------------|--|------------------------------|
| Sample Date | NYSDEC Groundwater Guidance or Standard Value (Note 1) | 9/3/2014 |
| Sample ID | | VPB154-GW-090314- 918-920 |
| Sample Interval | | 918 - 920 ft |
| Sample type code | | N |
| VOC 8260C (ug/L) | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 25 UJ |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 25 UJ |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | < 25 UJ |
| 1,1,2-TRICHLOROETHANE | 1 | < 25 UJ |
| 1,1-DICHLOROETHANE | 5 | < 25 UJ |
| 1,1-DICHLOROETHENE | 5 | < 25 UJ |
| 1,2,4-TRICHLOROBENZENE | 5 | < 25 UJ |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 38 UJ |
| 1,2-DIBROMOETHANE | NL | < 25 UJ |
| 1,2-DICHLOROBENZENE | 3 | < 25 UJ |
| 1,2-DICHLOROETHANE | 5 | < 25 UJ |
| 1,2-DICHLOROETHENE, TOTAL | 5 | < 50 UJ |
| 1,2-DICHLOROPROPANE | 1 | < 25 UJ |
| 1,3-DICHLOROBENZENE | 3 | < 25 UJ |
| 1,4-DICHLOROBENZENE | 3 | < 25 UJ |
| 2-BUTANONE | 50 | < 120 UJ |
| 2-HEXANONE | 50 | < 120 UJ |
| 4-METHYL-2-PENTANONE | NL | < 120 UJ |
| ACETONE | 50 | < 120 UJ |
| BENZENE | 1 | < 25 UJ |
| BROMODICHLOROMETHANE | 50 | < 25 UJ |
| BROMOFORM | 50 | < 25 UJ |
| BROMOMETHANE | 5 | < 50 UJ |
| CARBON DISULFIDE | 60 | < 25 UJ |
| CARBON TETRACHLORIDE | 5 | < 25 UJ |
| CHLOROBENZENE | 5 | < 25 UJ |
| CHLOROETHANE | 5 | < 50 UJ |
| CHLOROFORM | 7 | < 25 UJ |
| CHLOROMETHANE | 5 | < 50 UJ |
| CIS-1,2-DICHLOROETHENE | 5 | < 25 UJ |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 25 UJ |
| CYCLOHEXANE | NL | < 25 UJ |
| DIBROMOCHLOROMETHANE | 5 | < 25 UJ |
| DICHLORODIFLUOROMETHANE | 5 | < 50 UJ |
| ETHYLBENZENE | 5 | < 25 UJ |
| ISOPROPYLBENZENE | 5 | < 25 UJ |
| M- AND P-XYLENE | NL | < 50 UJ |
| METHYL ACETATE | NL | < 38 UJ |
| METHYL CYCLOHEXANE | NL | < 25 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 25 UJ |
| METHYLENE CHLORIDE | 5 | < 120 UJ |
| O-XYLENE | NL | < 25 UJ |
| STYRENE | 5 | < 25 UJ |
| TETRACHLOROETHENE | 5 | < 25 UJ |
| TOLUENE | 5 | < 25 UJ |
| TRANS-1,2-DICHLOROETHENE | 5 | < 25 UJ |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 25 UJ |
| TRICHLOROETHENE | 5 | < 25 UJ |
| TRICHLOROFLUOROMETHANE | 5 | < 50 UJ |
| VINYL CHLORIDE | 2 | < 50 UJ |
| XYLENES, TOTAL | 5 | < 75 UJ |

Notes:

1 New York State Department of Environmental Conservation Division of Water Technical and Operation Guidance series (6 NYCRR 700-706, Part 703.5 summarized in TOGS 1.1.1)

Ambient water quality standards and groundwater effluent limitations, class GA; NL = Not Listed

Bold = Detected; ***Bold and Italic***=Detection limit exceeds NYS Groundwater Standards or guidance value

Yellow highlighted values exceed Groundwater Standards or guidance value

Sample type codes: N - normal environmental sample, FD - field duplicate

U = Nondetected result. The analyte was analyzed for, but was not detected above the reported sample quantitation limit.

UJ = The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte.

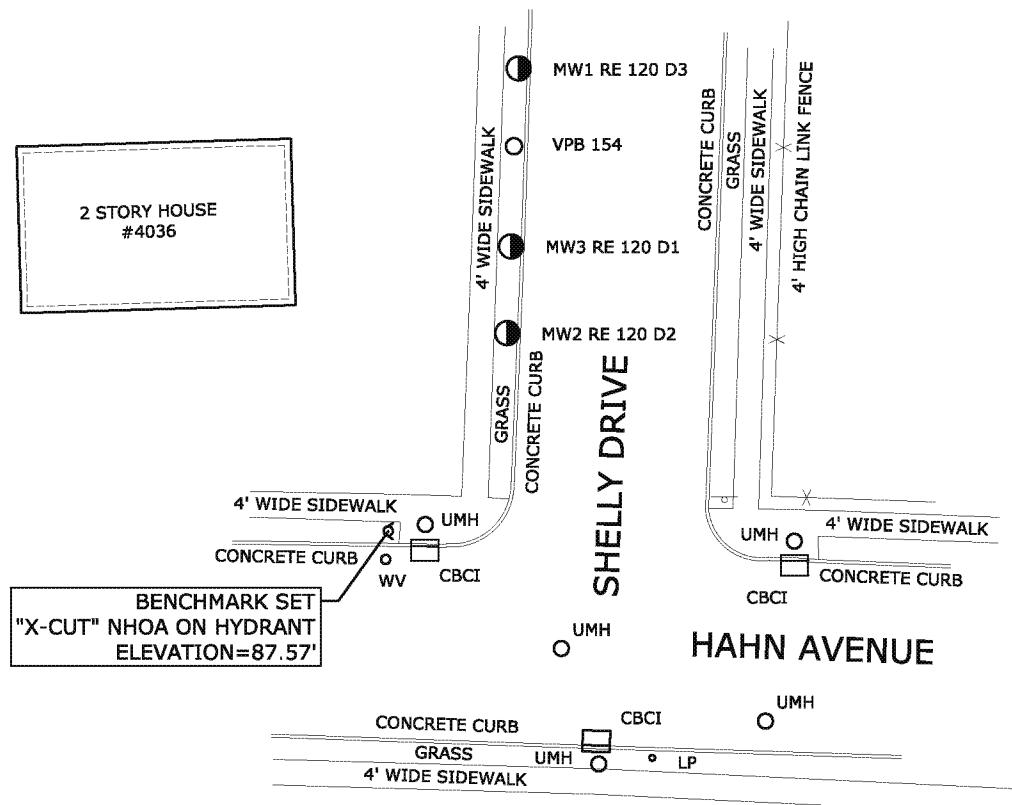
J = The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

Section 6

Survey

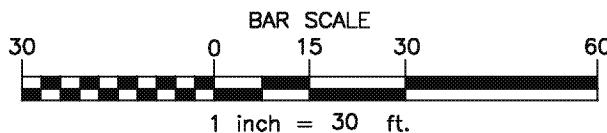
UNAUTHORIZED ALTERATION OR ADDITION TO
THIS DOCUMENT IS A VIOLATION OF SECTION
7209 SUBDIVISION 2 OF THE NEW YORK STATE
EDUCATION LAW.

| Description | Northing | Easting | Latitude | Longitude | Ground | Rim | PVC |
|---------------|-----------|------------|--------------|--------------|--------|-------|-------|
| MW1 RE 120 D3 | 204618.12 | 1125061.88 | N40-43-37.87 | W73-29-31.05 | 86.14 | 86.14 | 85.70 |
| VPB 154 | 204605.97 | 1125061.14 | N40-43-37.75 | W73-29-31.06 | 85.86 | NA | NA |
| MW3 RE 120 D1 | 204590.37 | 1125060.70 | N40-43-37.59 | W73-29-31.07 | 86.06 | 86.06 | 85.58 |
| MW2 RE 120 D2 | 204576.78 | 1125060.08 | N40-43-37.46 | W73-29-31.08 | 86.03 | 86.03 | 85.54 |



Map Notes

- Information shown hereon was compiled from an actual field survey conducted from December 9, 2014.
- North orientation is Grid North based on the New York State Plane Coordinate System, Long Island Zone, NAD 83 as obtained from GPS observations.
- Vertical datum shown hereon is NAVD 88 as obtained from GPS observations.



Legend

| | | |
|--------------------------|---------|-------------------------|
| <input type="checkbox"/> | CBCI | Catch Basin Curb Inlet |
| ○ | GV | Gas Valve |
| ○ | LP | Light Post |
| ● | MW1 | Monitoring Well |
| — | Sign | |
| ○ | UMH | Unknown Manhole |
| ○ | VPB 149 | Vertical Profile Boring |
| ○ | WV | Water Valve |

DWG NO. 14-644

| Date | RECORD OF WORK | Appr. | VERICAL PROFILE BORING 154 SURVEY LOCATION SHELLY DRIVE | |
|---------------|-------------------|-------|---|-------------------------|
| | | | | |
| | | | | |
| | | | | |
| | | | | |
| | | | | |
| Drafter: LMK | Checker: JFC | | TOWN OF BETHPAGE | NASSAU COUNTY, NEW YORK |
| Appr. by: JFC | Proj. No. 14.4121 | | C.T. MALE ASSOCIATES Engineering, Surveying, Architecture & Landscape Architecture, D.P.C. | |
| | | | 50 CENTURY HILL DRIVE, LATHAM, NY 12110 518.786.7400 * FAX 518.786.7299 | |
| | | | SCALE: 1"=30' | DATE: DEC. 09, 2014 |

